

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG059004.D
 Acq On : 12 Sep 2023 6:29
 Operator : MA/JU
 Sample : 04235-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EZYJ9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG059004.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.638	699	706	712	rBV	134915	229491	1.71%	0.125%
2	8.325	816	823	830	rVV	417390	733569	5.48%	0.400%
3	8.913	915	923	932	rBV7	47902	131864	0.98%	0.072%
4	9.007	932	939	945	rBV7	72490	171396	1.28%	0.093%
5	9.160	958	965	975	rVB8	79295	208157	1.55%	0.113%
6	9.383	993	1003	1012	rBV8	90735	306338	2.29%	0.167%
7	9.483	1012	1020	1024	rBV	368250	653772	4.88%	0.356%
8	9.741	1057	1064	1072	rVB6	99285	224834	1.68%	0.123%
9	9.900	1085	1091	1095	rBV3	145927	264644	1.98%	0.144%
10	10.035	1110	1114	1124	rVB4	173526	336234	2.51%	0.183%
11	10.135	1126	1131	1135	rBV3	94246	163699	1.22%	0.089%
12	10.211	1138	1144	1153	rVB4	187241	448592	3.35%	0.244%
13	10.305	1153	1160	1171	rVB5	323370	938141	7.00%	0.511%
14	10.405	1171	1177	1182	rBV2	215766	399142	2.98%	0.217%
15	10.493	1185	1192	1196	rBV3	221554	381820	2.85%	0.208%
16	10.593	1203	1209	1214	rVB2	236128	423750	3.16%	0.231%
17	10.799	1239	1244	1252	rVB7	116279	310953	2.32%	0.169%
18	10.940	1265	1268	1273	rVV5	172621	346799	2.59%	0.189%
19	10.993	1273	1277	1281	rVV4	207842	438027	3.27%	0.239%
20	11.046	1281	1286	1291	rVV	2339137	3738178	27.91%	2.037%
21	11.093	1291	1294	1298	rVV4	203096	341387	2.55%	0.186%
22	11.169	1298	1307	1312	rVV	738150	1745546	13.03%	0.951%
23	11.228	1312	1317	1324	rVV	1486679	2442621	18.23%	1.331%
24	11.316	1324	1332	1344	rVB8	312730	1136659	8.49%	0.619%
25	11.457	1354	1356	1365	rVB8	108086	223902	1.67%	0.122%
26	11.592	1365	1379	1382	rBV8	307946	971202	7.25%	0.529%
27	11.727	1399	1402	1405	rBV4	122528	176613	1.32%	0.096%
28	11.804	1409	1415	1426	rVB4	839928	2080484	15.53%	1.134%
29	11.909	1428	1433	1437	rBV	379400	669200	5.00%	0.365%
30	11.986	1442	1446	1453	rVB	540882	939714	7.01%	0.512%
31	12.086	1458	1463	1468	rVV	2358204	3395876	25.35%	1.850%
32	12.144	1469	1473	1481	rVB5	279235	632199	4.72%	0.344%
33	12.303	1495	1500	1505	rBV4	208363	380868	2.84%	0.208%
34	12.362	1505	1510	1516	rVB4	435702	800430	5.98%	0.436%
35	12.426	1516	1521	1525	rBV4	355754	766116	5.72%	0.417%
36	12.485	1525	1531	1539	rVB	6570323	9684612	72.30%	5.277%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG059004.D
 Acq On : 12 Sep 2023 6:29
 Operator : MA/JU
 Sample : 04235-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EZYJ9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Title : SVOA CALIBRATION

37	12.556	1539	1543	1549	rBV7	282317	661293	4.94%	0.360%
38	12.608	1549	1552	1555	rBV2	207116	314056	2.34%	0.171%
39	12.691	1560	1566	1574	rVB	952877	1901780	14.20%	1.036%
40	12.796	1579	1584	1588	rBV	330407	517390	3.86%	0.282%
41	12.896	1596	1601	1607	rBV7	409581	1094700	8.17%	0.597%
42	13.014	1616	1621	1624	rBV	360204	600180	4.48%	0.327%
43	13.102	1631	1636	1639	rBV2	751246	1218559	9.10%	0.664%
44	13.137	1639	1642	1645	rVV2	569119	821173	6.13%	0.447%
45	13.190	1645	1651	1659	rVB3	893995	2571371	19.20%	1.401%
46	13.278	1661	1666	1673	rBV	1201401	1854170	13.84%	1.010%
47	13.331	1673	1675	1677	rBV2	123072	141071	1.05%	0.077%
48	13.366	1677	1681	1684	rVV	949219	1278528	9.54%	0.697%
49	13.413	1684	1689	1697	rVB	3106483	4423481	33.02%	2.410%
50	13.707	1733	1739	1744	rBV	8863979	13395799	100.00%	7.300%
51	13.760	1744	1748	1749	rVV2	370674	561873	4.19%	0.306%
52	13.784	1750	1752	1758	rVV4	590395	1011345	7.55%	0.551%
53	13.860	1762	1765	1768	rVB3	292488	343008	2.56%	0.187%
54	13.942	1775	1779	1785	rBV7	208874	476246	3.56%	0.260%
55	14.101	1801	1806	1807	rBV2	551359	762976	5.70%	0.416%
56	14.224	1823	1827	1830	rBV3	308853	467479	3.49%	0.255%
57	14.265	1830	1834	1839	rVB3	1252047	2023796	15.11%	1.103%
58	14.348	1839	1848	1852	rBV3	4238758	8051871	60.11%	4.388%
59	14.471	1865	1869	1872	rBV	683100	780141	5.82%	0.425%
60	14.653	1895	1900	1905	rVB5	840180	1356695	10.13%	0.739%
61	14.771	1915	1920	1927	rVB	8343048	12067148	90.08%	6.576%
62	14.853	1928	1934	1937	rBV5	476432	970926	7.25%	0.529%
63	14.947	1946	1950	1954	rBV	662377	1061221	7.92%	0.578%
64	15.017	1959	1962	1964	rBV3	219431	297383	2.22%	0.162%
65	15.205	1990	1994	1996	rBV	483643	731422	5.46%	0.399%
66	15.258	2000	2003	2007	rVB	809635	994079	7.42%	0.542%
67	15.317	2009	2013	2018	rVB4	839538	1190375	8.89%	0.649%
68	15.376	2019	2023	2026	rBV	1532036	2086825	15.58%	1.137%
69	15.446	2033	2035	2041	rVB	598406	821256	6.13%	0.448%
70	15.540	2048	2051	2056	rVV3	343031	545971	4.08%	0.298%
71	15.599	2057	2061	2068	rVB5	691720	1328829	9.92%	0.724%
72	15.717	2074	2081	2085	rBV	6722778	8971522	66.97%	4.889%
73	15.769	2087	2090	2094	rVV	478744	626493	4.68%	0.341%
74	15.828	2098	2100	2104	rVB3	306013	347271	2.59%	0.189%
75	15.934	2114	2118	2127	rVB4	574163	1381694	10.31%	0.753%
76	16.016	2128	2132	2138	rVB5	297307	571881	4.27%	0.312%

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG059004.D
 Acq On : 12 Sep 2023 6:29
 Operator : MA/JU
 Sample : 04235-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EZYJ9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Title : SVOA CALIBRATION

77	16.110	2142	2148	2153	rBV	2359953	3966645	29.61%	2.161%
78	16.157	2153	2156	2159	rBV6	189902	254208	1.90%	0.139%
79	16.269	2172	2175	2178	rBV3	442409	623905	4.66%	0.340%
80	16.339	2183	2187	2195	rVB2	552696	969264	7.24%	0.528%
81	16.580	2222	2228	2235	rBV	5183113	9807634	73.21%	5.344%
82	16.845	2270	2273	2276	rVB2	259180	279438	2.09%	0.152%
83	17.086	2311	2314	2321	rVB2	452580	624851	4.66%	0.340%
84	17.150	2322	2325	2328	rVV3	348359	432290	3.23%	0.236%
85	17.203	2331	2334	2343	rVV5	392227	905155	6.76%	0.493%
86	17.373	2358	2363	2366	rBV	2967703	3533198	26.38%	1.925%
87	17.415	2366	2370	2375	rVB	1787308	2655371	19.82%	1.447%
88	17.697	2415	2418	2423	rVB	858789	1132801	8.46%	0.617%
89	17.796	2432	2435	2440	rVB2	580963	708520	5.29%	0.386%
90	17.996	2466	2469	2470	rBV	490066	477377	3.56%	0.260%
91	18.114	2485	2489	2494	rBV	3404308	4502027	33.61%	2.453%
92	18.560	2561	2565	2569	rBV3	939712	1433777	10.70%	0.781%
93	18.613	2572	2574	2577	rVB2	644704	770941	5.76%	0.420%
94	18.801	2602	2606	2611	rBV	3734893	4873186	36.38%	2.655%
95	19.430	2709	2713	2717	rBV	3956506	4720194	35.24%	2.572%
96	19.853	2782	2785	2790	rVB	1419551	1951424	14.57%	1.063%
97	20.000	2806	2810	2815	rBV2	3826048	5128281	38.28%	2.794%
98	20.540	2898	2902	2906	rBV3	2796324	4463966	33.32%	2.432%
99	20.711	2928	2931	2937	rBV	3266801	5176962	38.65%	2.821%
100	21.780	3109	3113	3120	rBV3	3004629	6262325	46.75%	3.412%

Sum of corrected areas: 183513846

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

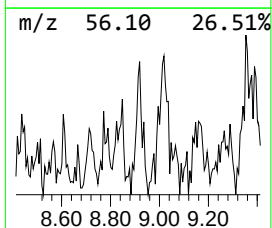
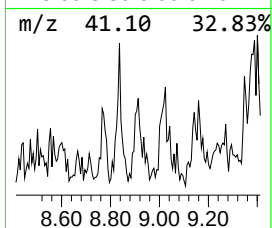
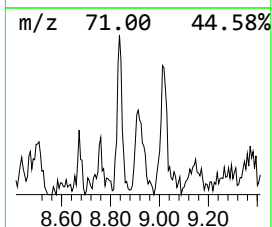
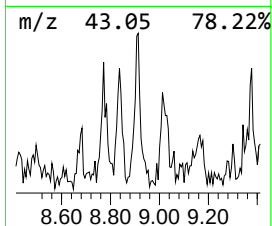
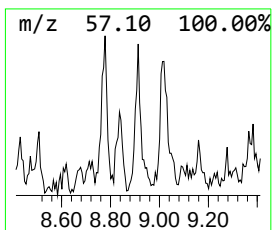
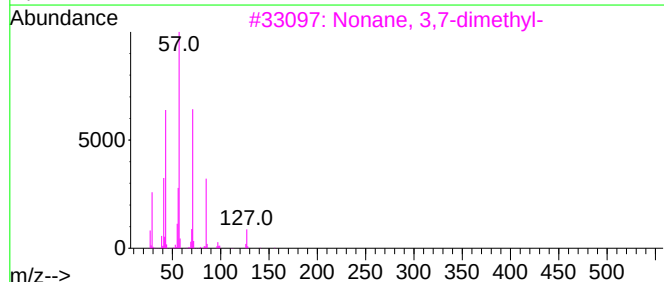
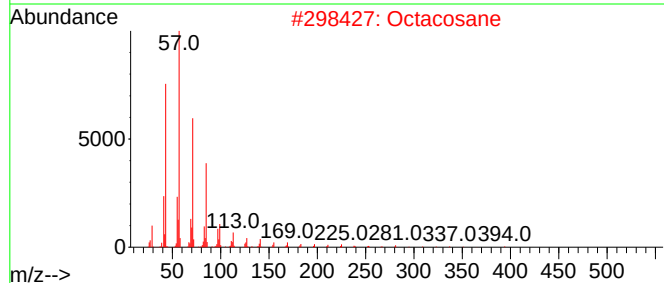
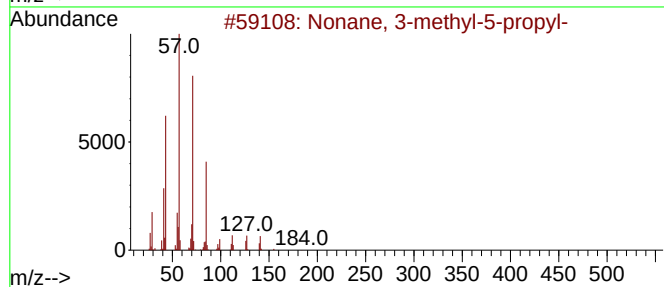
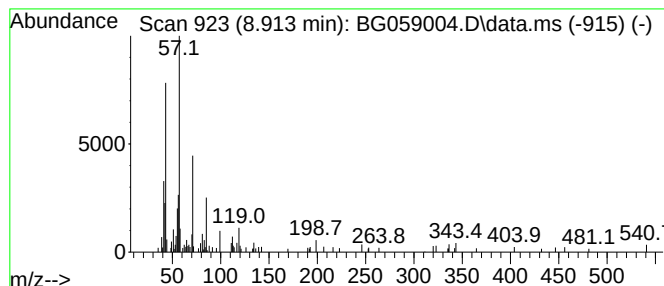
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 77

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.913	3.60 ng/ul	131864	1,4-Dichlorobenzene-d4			8.325
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-5-propyl-		184	C13H28	031081-18-2	53
2	Octacosane		394	C28H58	000630-02-4	53
3	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	53
4	Undecane, 3-methyl-		170	C12H26	001002-43-3	53
5	Decane, 6-ethyl-2-methyl-		184	C13H28	062108-21-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

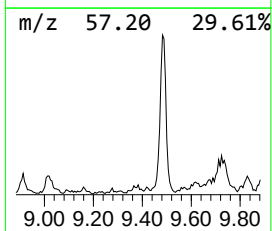
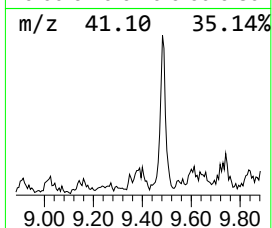
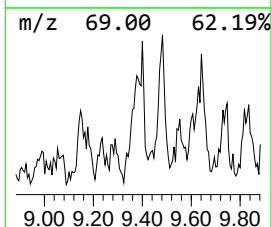
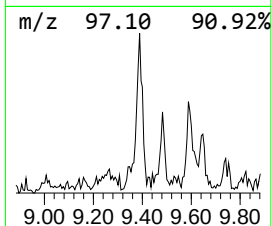
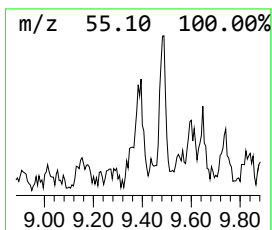
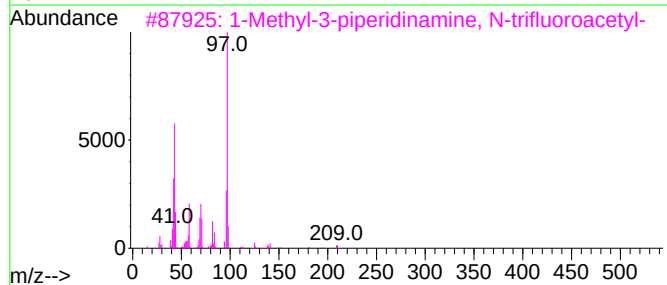
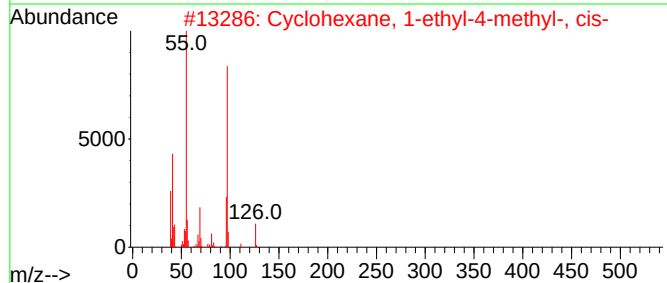
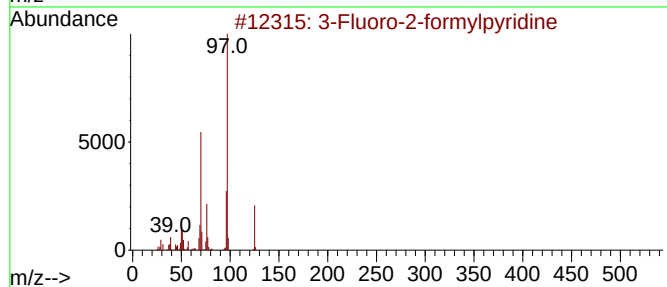
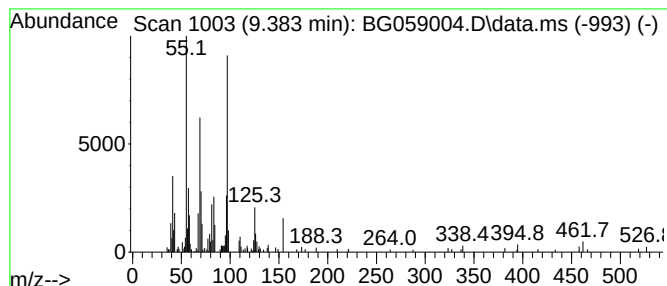
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 3-Fluoro-2-formylpyridine Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.383	8.35 ng/ul	306338	1,4-Dichlorobenzene-d4	8.325

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Fluoro-2-formylpyridine	125	C6H4FN0	031224-43-8	72
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	70
3			1-Methyl-3-piperidinamine, N-tri...	210	C8H13F3N2O	1344377-30-5	58
4			Oxalic acid, butyl cyclohexylmet...	242	C13H22O4	1000309-68-2	58
5			Oxalic acid, cyclohexylmethyl is...	270	C15H26O4	1000309-68-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

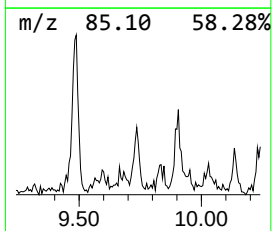
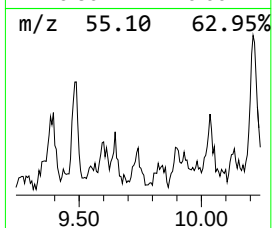
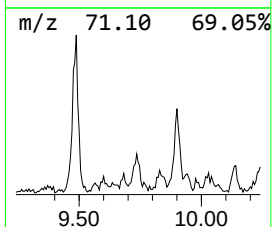
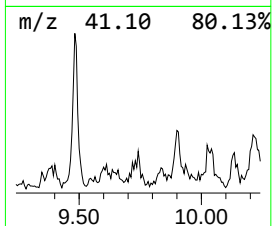
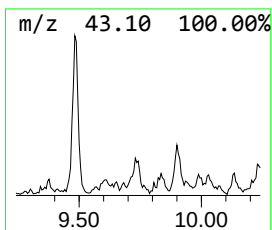
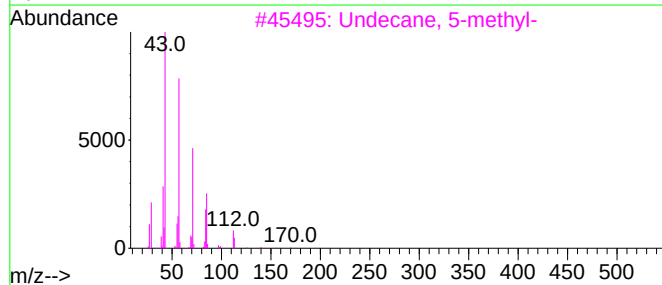
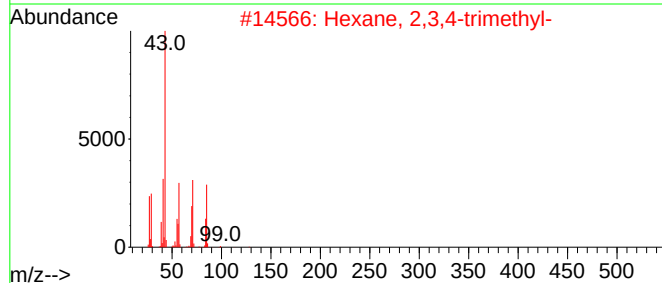
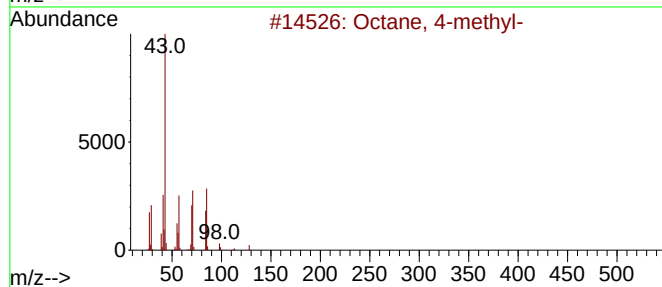
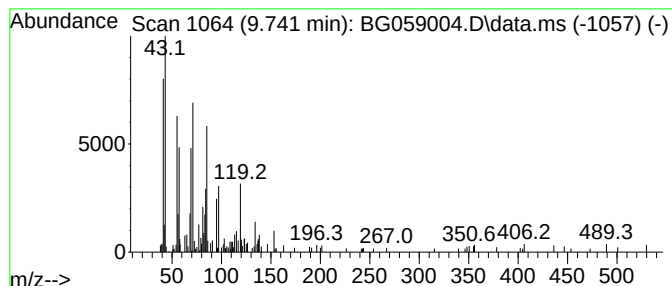
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 63

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.741	6.13 ng/ul	224834	1,4-Dichlorobenzene-d4	8.325

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-methyl-			128	C9H20	002216-34-4	43
2	Hexane, 2,3,4-trimethyl-			128	C9H20	000921-47-1	43
3	Undecane, 5-methyl-			170	C12H26	001632-70-8	43
4	Pentadecane, 5-methyl-			226	C16H34	025117-33-3	38
5	Octadecane, 5,14-dibutyl-			366	C26H54	055282-13-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

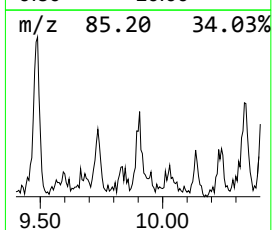
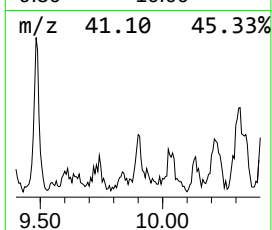
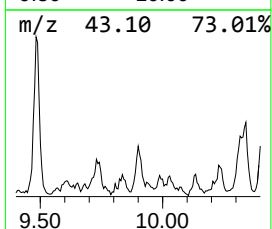
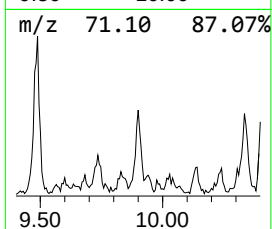
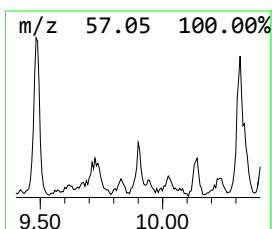
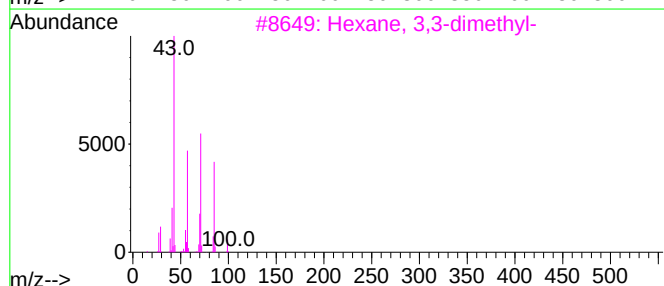
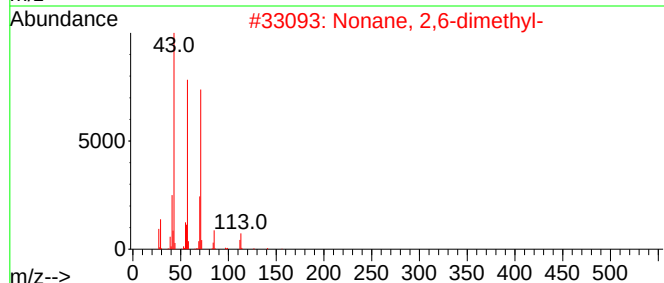
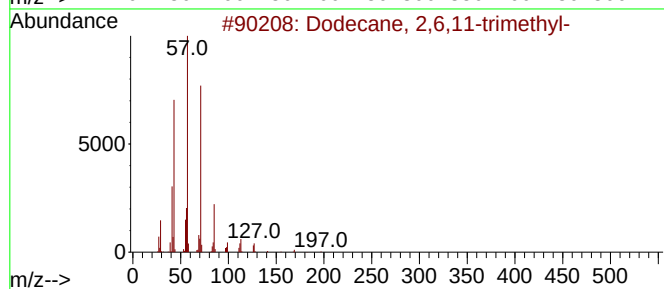
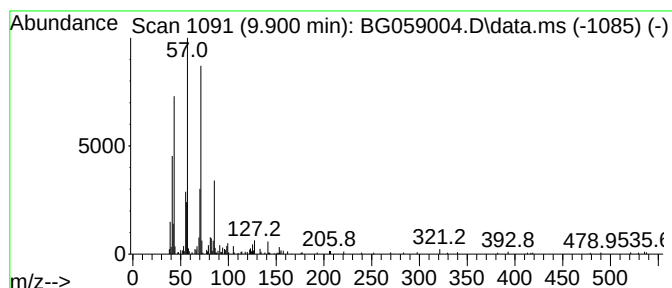
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 79

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.900	3.03 ng/ul	264644	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
2	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	64
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	Decane, 4-methyl-	156	C11H24	002847-72-5	53
5	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

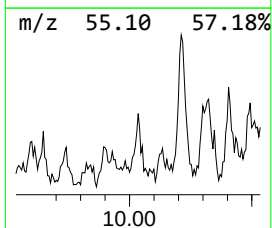
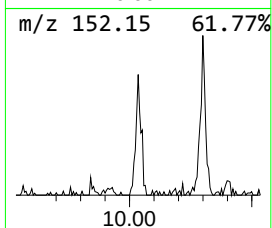
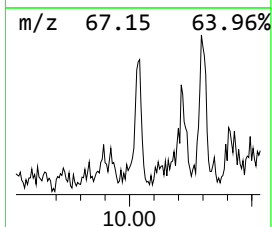
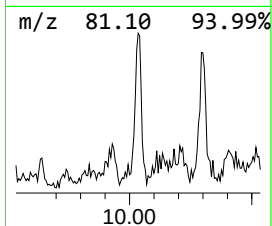
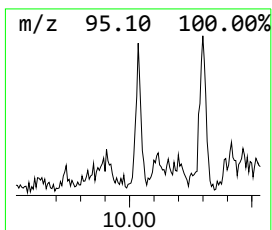
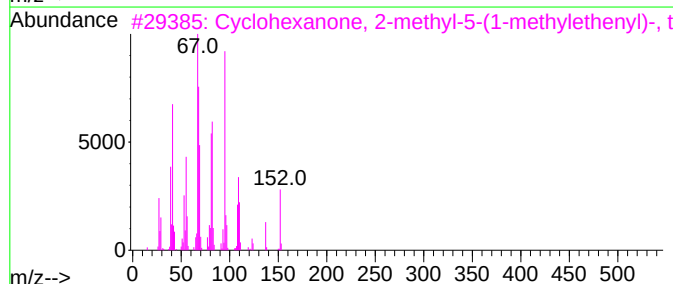
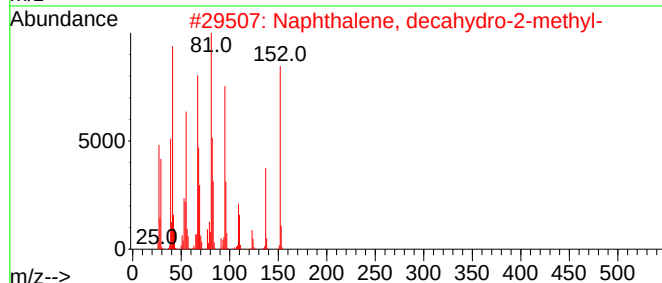
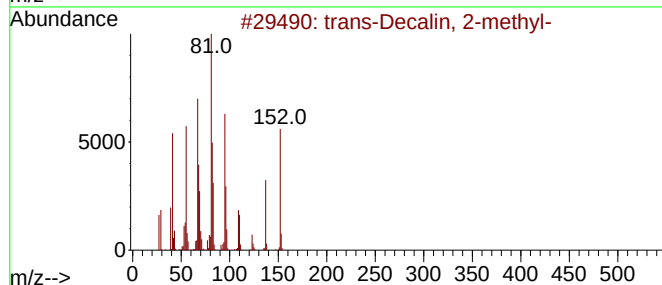
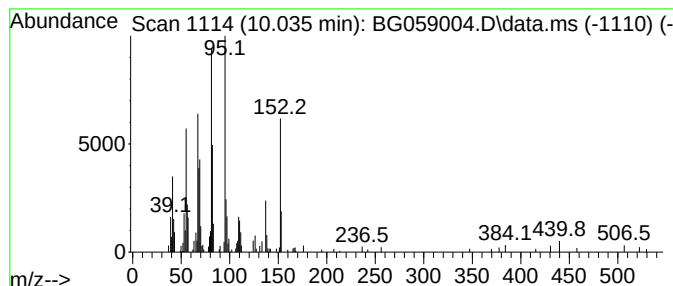
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 trans-Decalin, 2-methyl- Concentration Rank 75

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.035	3.85 ng/ul	336234	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	90
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	89
3			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	64
4			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	007764-50-3	64
5			(+)-Dihydrocarvone	152	C10H16O	005524-05-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

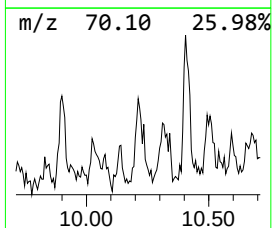
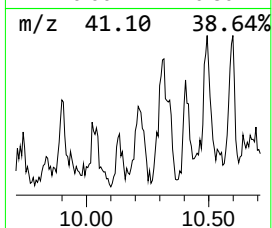
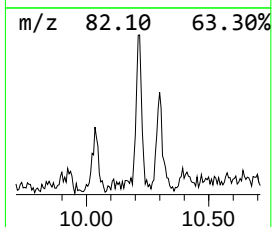
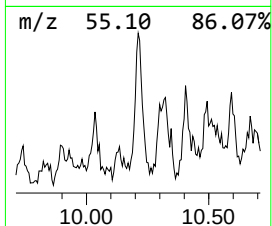
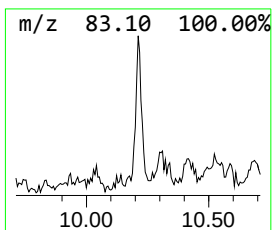
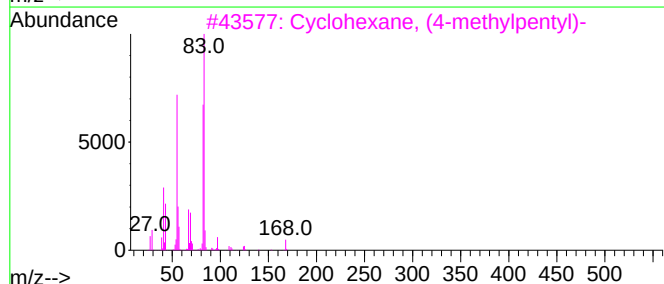
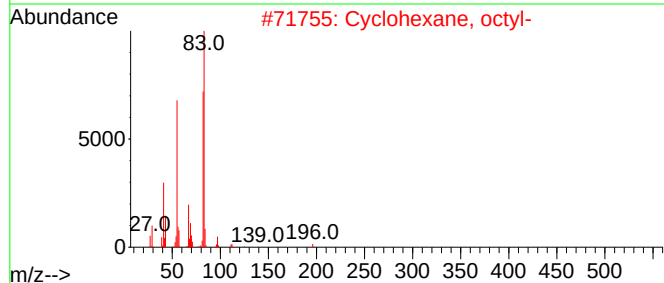
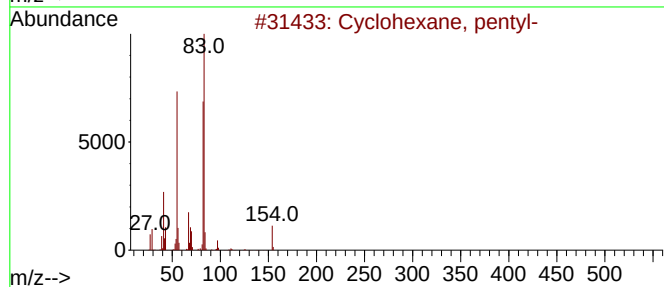
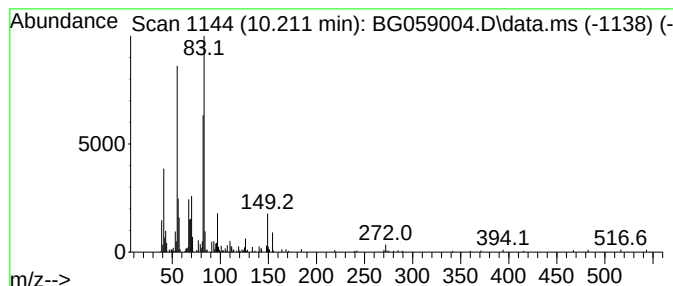
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Cyclic10.211 Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.211	5.14 ng/ul	448592	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, pentyl-	154	C11H22	004292-92-6	80
2	Cyclohexane, octyl-	196	C14H28	001795-15-9	59
3	Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	59
4	Cyclohexane, propyl-	126	C9H18	001678-92-8	59
5	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

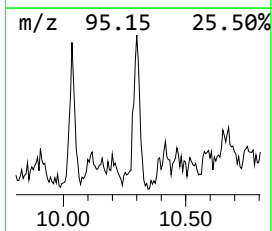
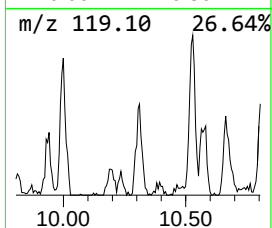
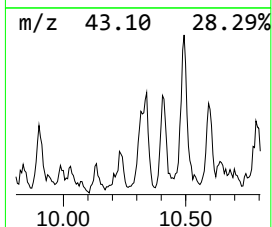
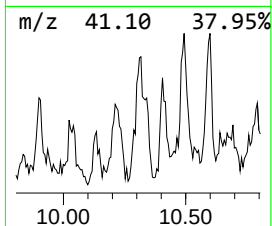
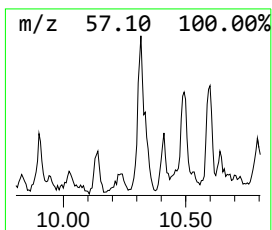
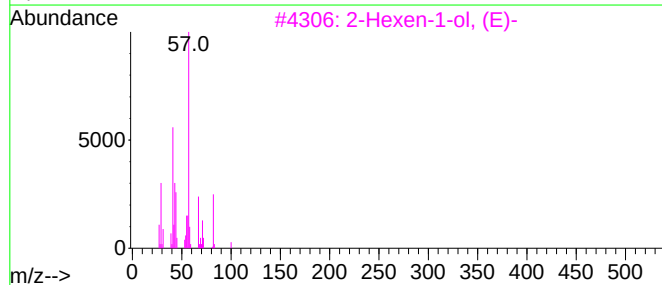
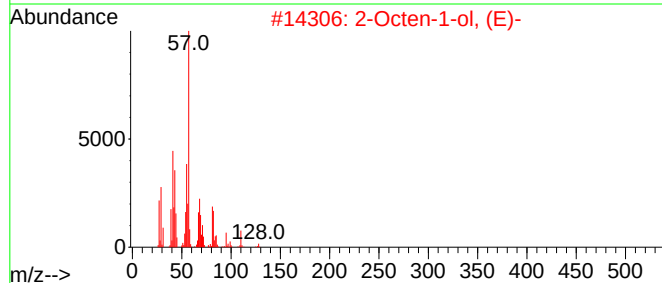
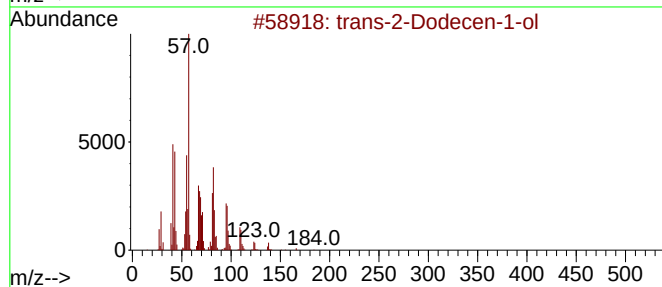
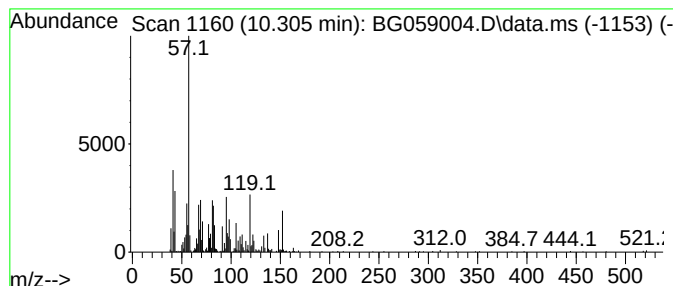
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-01 Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.305	10.75 ng/ul	938141	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-2-Dodecen-1-ol	184	C12H24O	069064-37-5	35
2			2-Octen-1-ol, (E)-	128	C8H16O	018409-17-1	35
3			2-Hexen-1-ol, (E)-	100	C6H12O	000928-95-0	27
4			2-Hexen-1-ol, (E)-	100	C6H12O	000928-95-0	27
5			2-Hexen-1-ol, propanoate, (E)-	156	C9H16O2	053398-80-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

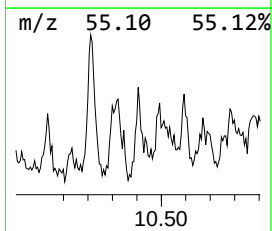
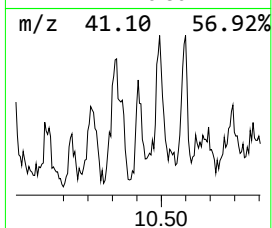
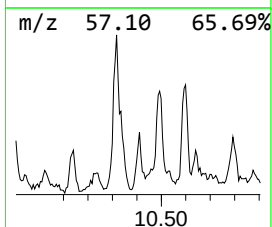
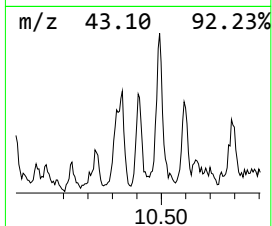
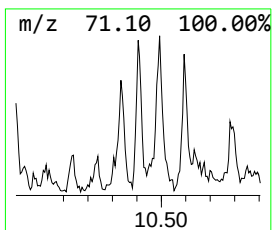
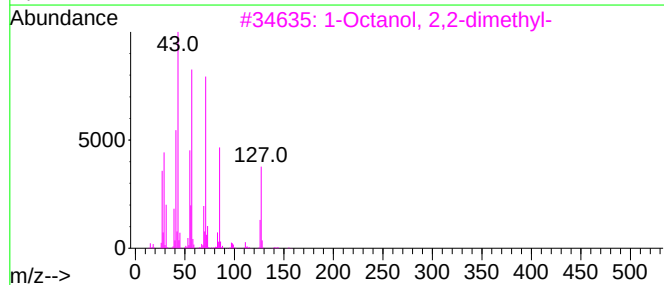
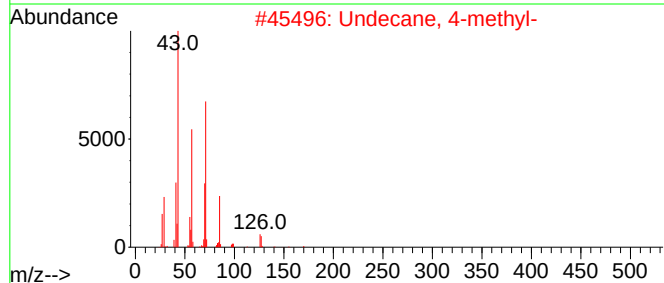
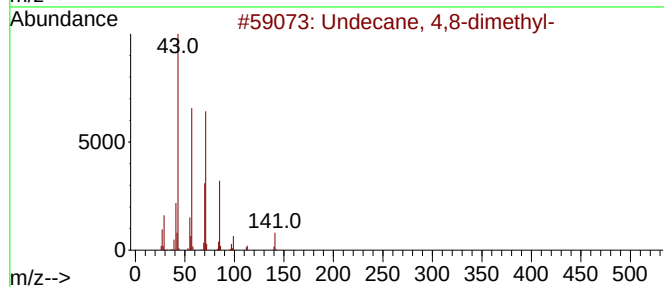
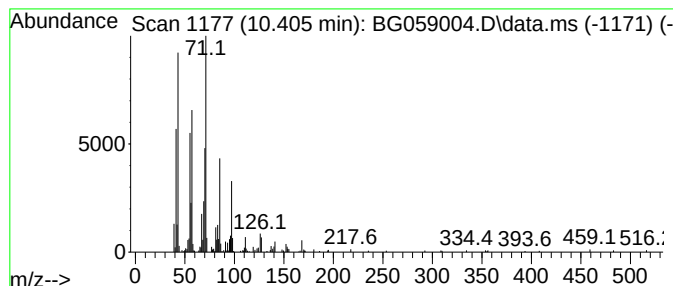
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 70

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.405	4.57 ng/ul	399142	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	59
2	Undecane, 4-methyl-	170	C12H26	002980-69-0	52
3	1-Octanol, 2,2-dimethyl-	158	C10H22O	002370-14-1	50
4	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	47
5	Propanoic acid, 2-methyl-, 2-eth...	200	C12H24O2	035061-61-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

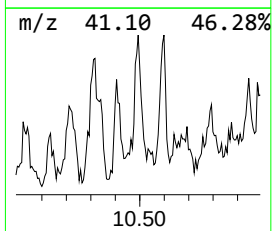
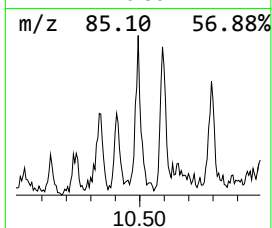
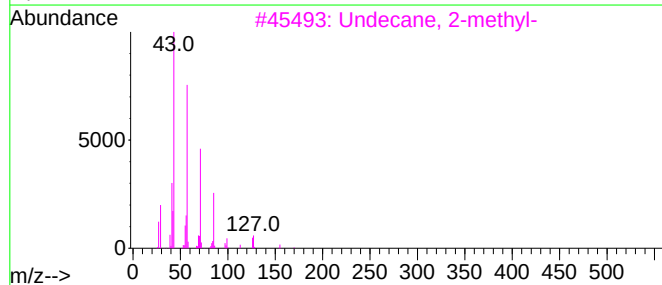
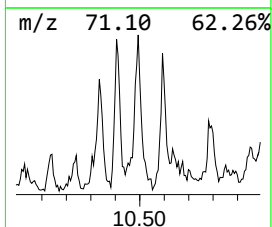
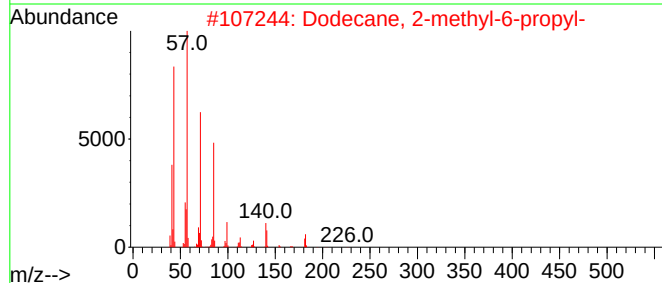
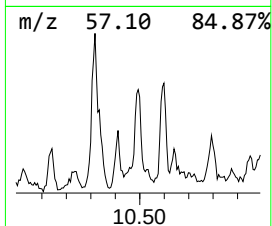
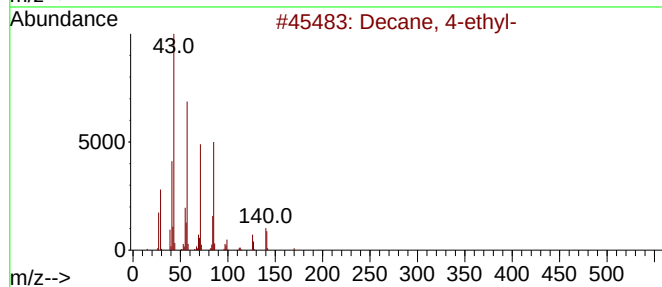
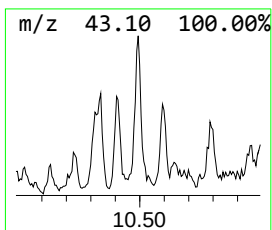
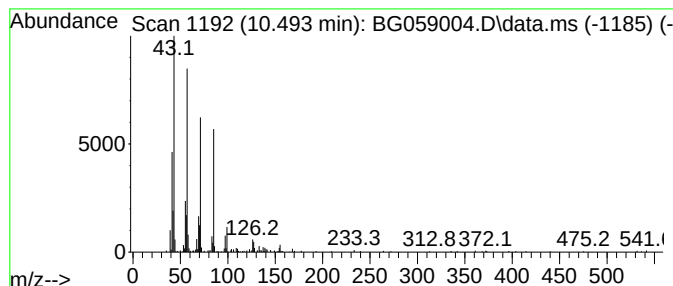
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 71

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.493	4.37 ng/ul	381820	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-ethyl-	170	C12H26	001636-44-8	64
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	59
3	Undecane, 2-methyl-	170	C12H26	007045-71-8	58
4	Hexadecane	226	C16H34	000544-76-3	53
5	Sulfurous acid, hexyl undecyl ester	320	C17H36O3S	1000309-13-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

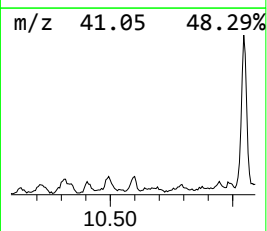
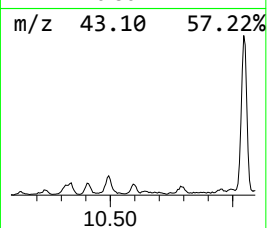
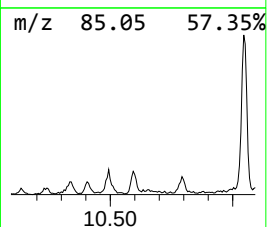
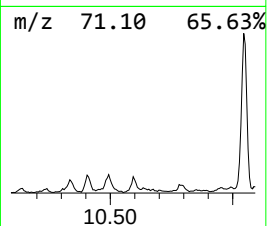
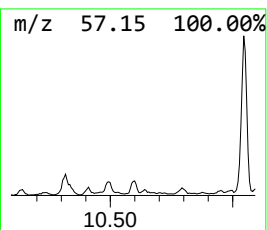
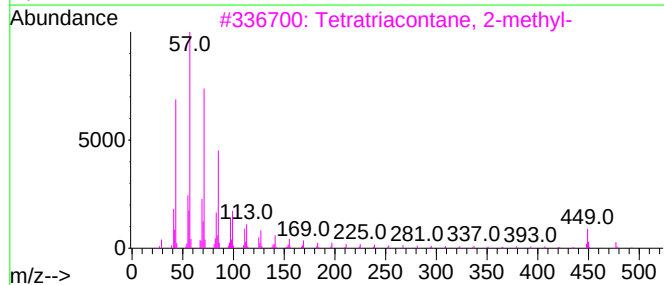
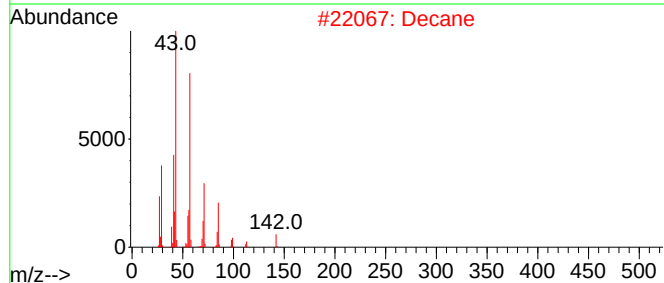
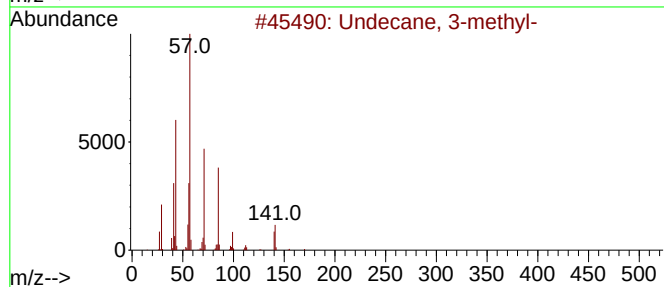
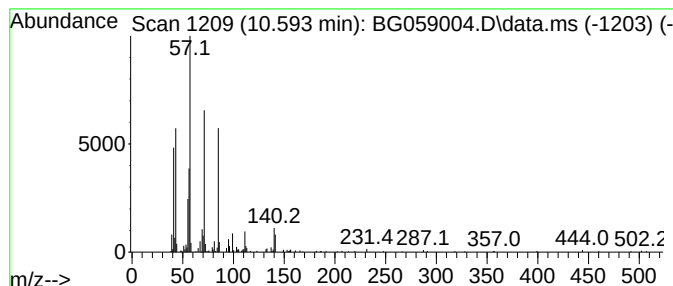
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 68

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.593	4.86 ng/ul	423750	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3-methyl-	170	C12H26	001002-43-3	59
2	Decane	142	C10H22	000124-18-5	53
3	Tetratriacontane, 2-methyl-	493	C35H72	014167-65-8	52
4	Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	50
5	Decane, 1-iodo-	268	C10H21I	002050-77-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

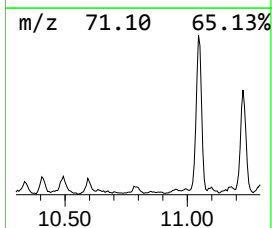
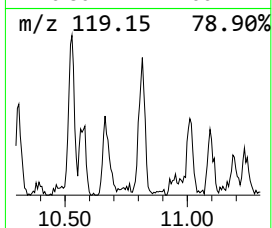
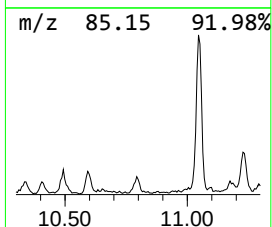
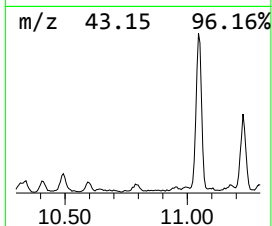
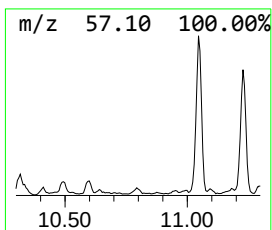
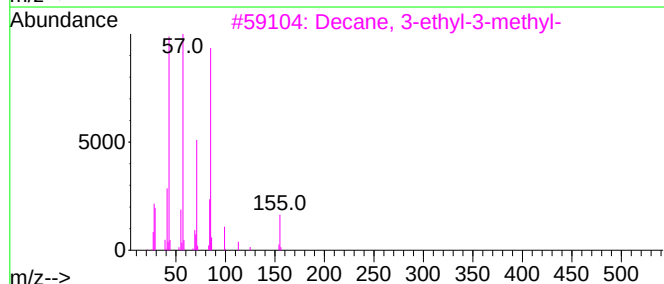
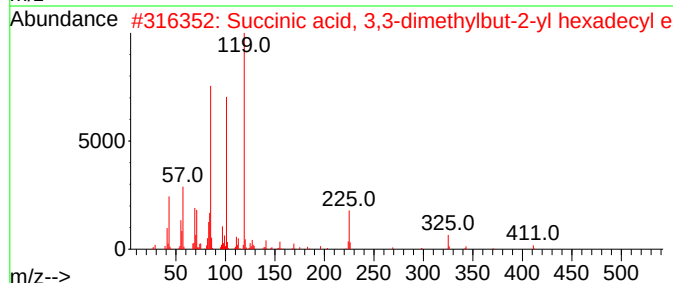
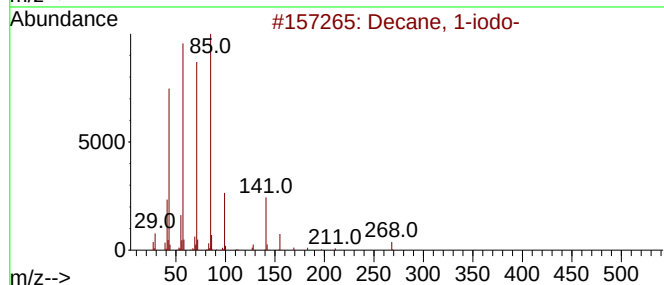
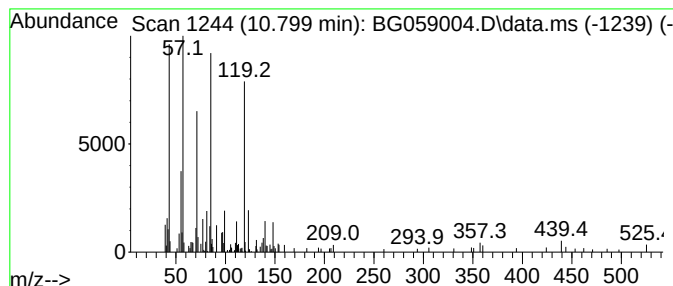
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown-02 Concentration Rank 78

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.799	3.56 ng/ul	310953	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 1-iodo-	268	C10H21I	002050-77-3	43
2			Succinic acid, 3,3-dimethylbut-2...	426	C26H50O4	1000349-52-0	38
3			Decane, 3-ethyl-3-methyl-	184	C13H28	017312-66-2	38
4			Sulfurous acid, 2-ethylhexyl iso...	278	C14H30O3S	1000309-19-0	35
5			Sulfurous acid, isohexyl pentyl ...	236	C11H24O3S	1000309-14-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

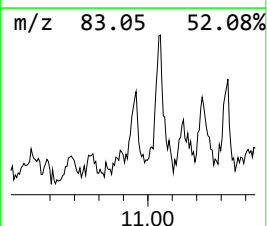
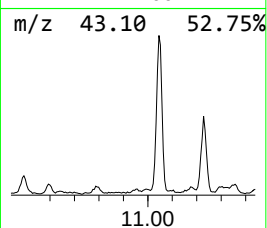
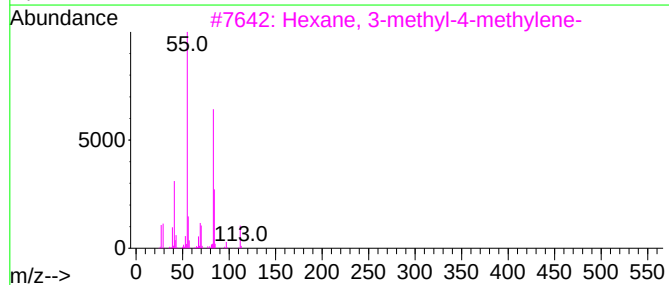
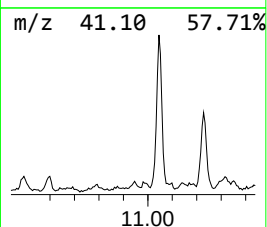
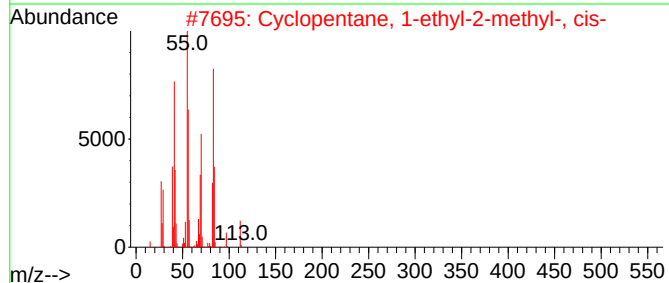
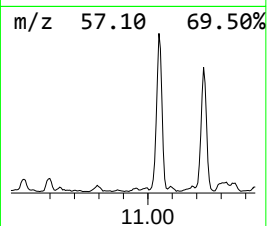
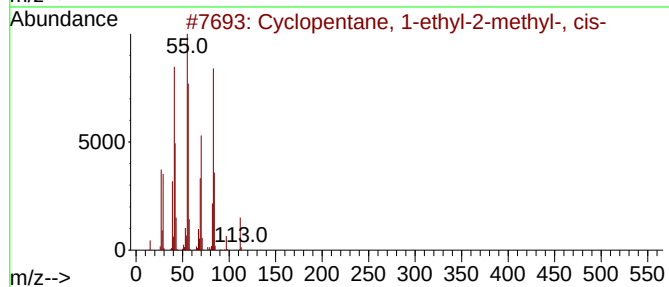
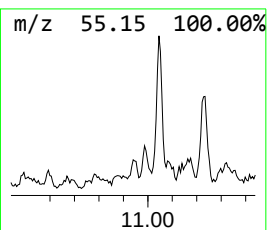
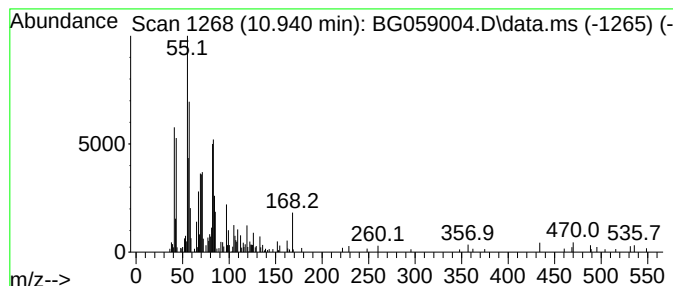
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Cyclic10.940 Concentration Rank 73

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.940	3.97 ng/ul	346799	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	49
2			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	47
3			Hexane, 3-methyl-4-methylene-	112	C8H16	003404-67-9	47
4			2-Dodecene, (E)-	168	C12H24	007206-13-5	46
5			2-Ethyl-1-hexanol, pentafluoropr...	276	C11H17F5O2	959012-82-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

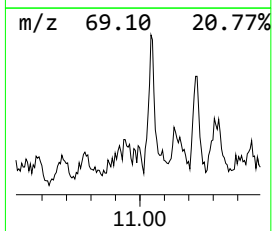
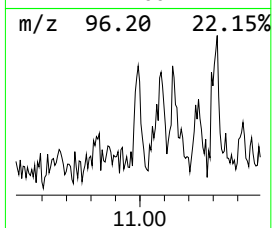
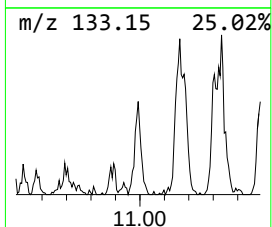
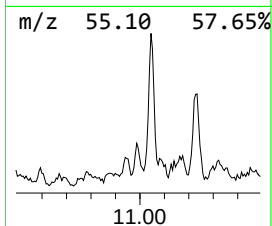
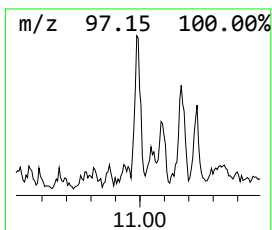
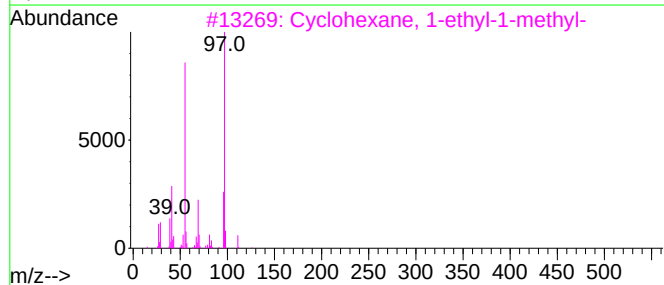
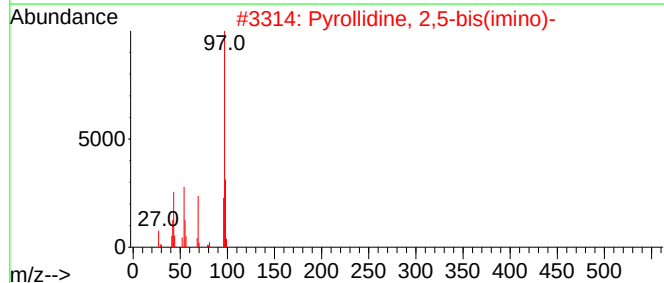
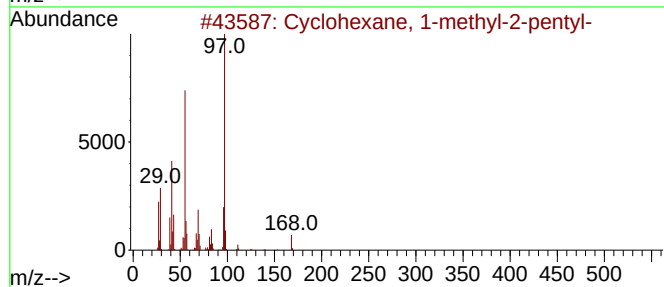
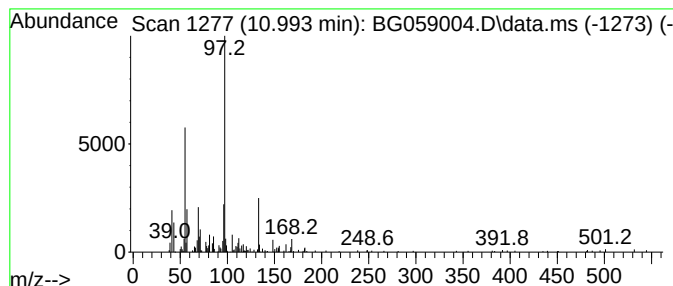
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Cyclic10.993 Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.993	5.02 ng/ul	438027	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	58
2			Pyrollidine, 2,5-bis(imino)-	97	C4H7N3	000503-84-4	53
3			Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	53
4			Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	53
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

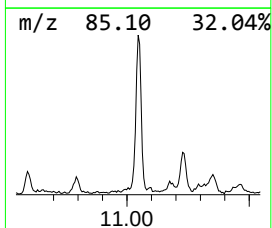
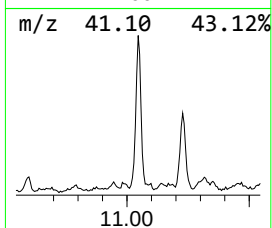
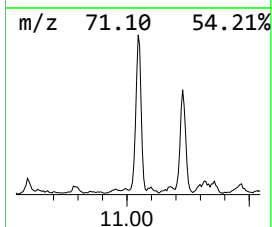
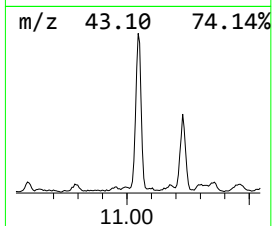
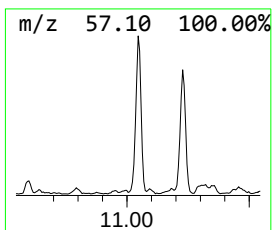
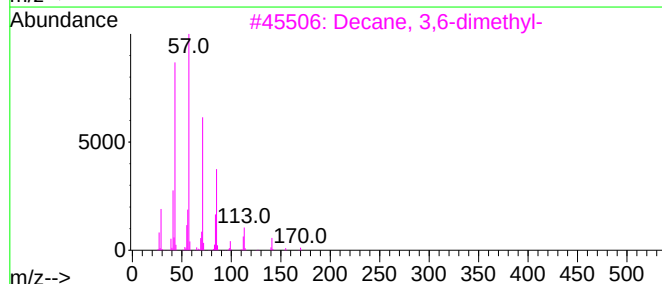
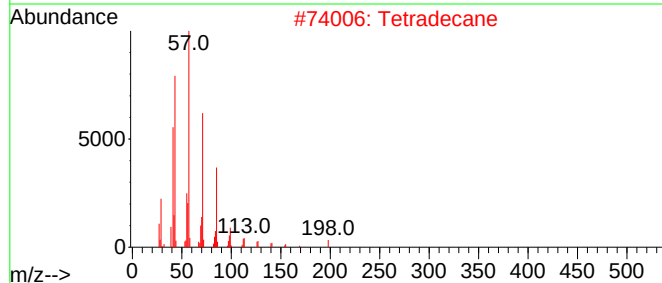
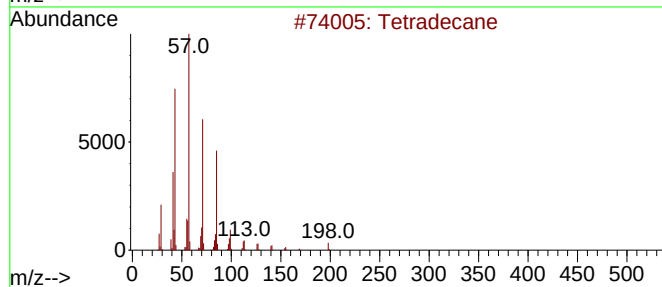
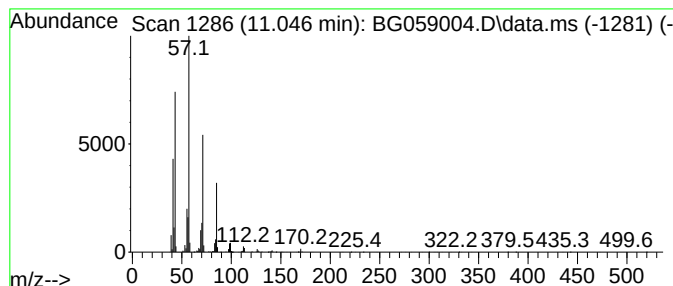
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.046	42.83 ng/ul	3738180	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	86
2	Tetradecane	198	C14H30	000629-59-4	83
3	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	80
4	Undecane, 5-methyl-	170	C12H26	001632-70-8	80
5	Hexadecane	226	C16H34	000544-76-3	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

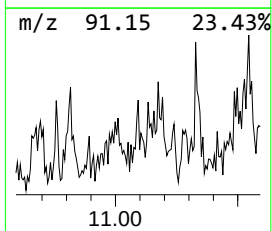
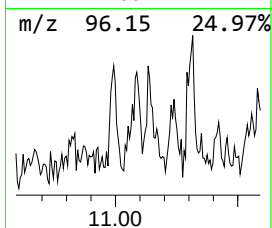
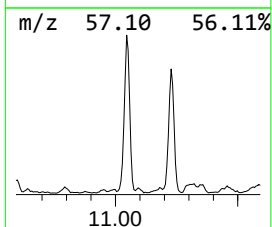
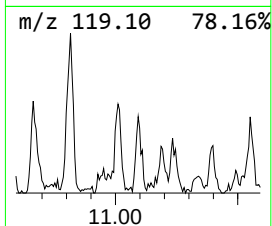
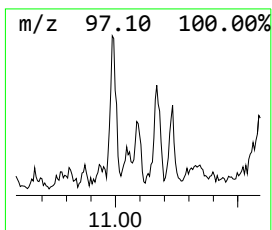
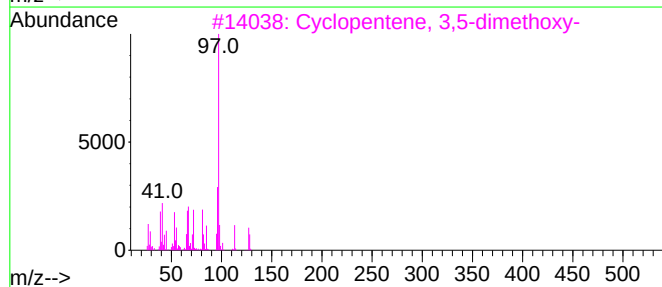
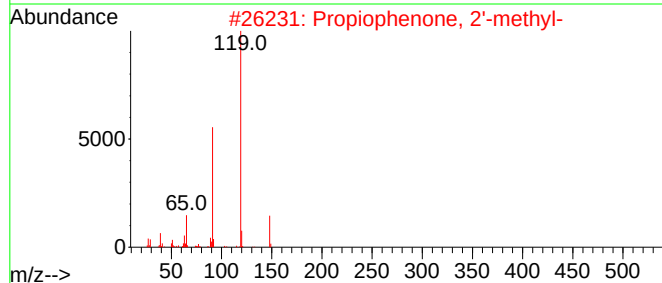
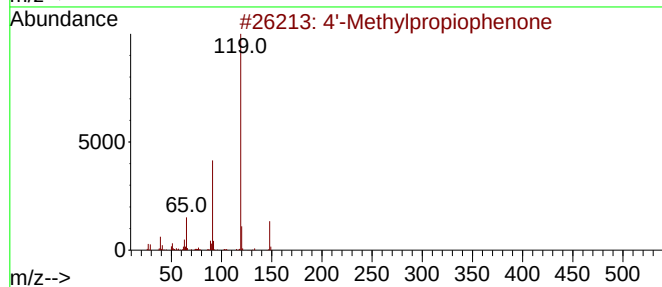
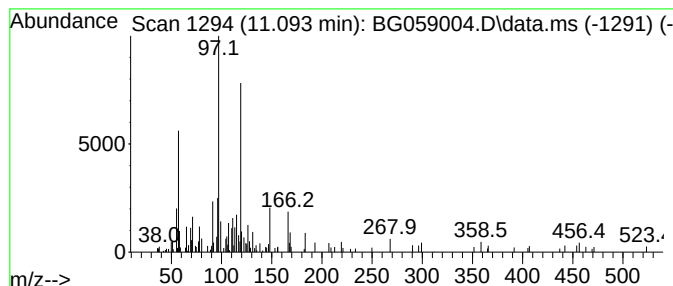
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown-03 Concentration Rank 74

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.093	3.91 ng/ul	341387	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4'-Methylpropiophenone	148	C10H12O	005337-93-9	38
2	Propiophenone, 2'-methyl-	148	C10H12O	002040-14-4	35
3	Cyclopentene, 3,5-dimethoxy-	128	C7H12O2	089897-05-2	35
4	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	35
5	Benzamide, 4-methyl-N-butyl-N-et...	219	C14H21NO	1000415-91-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

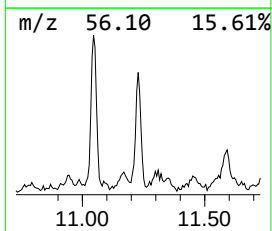
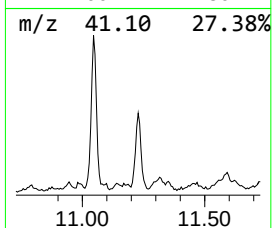
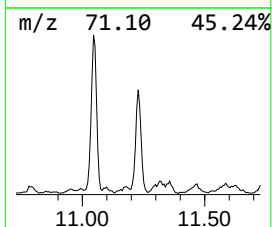
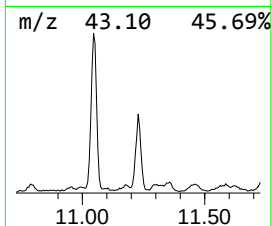
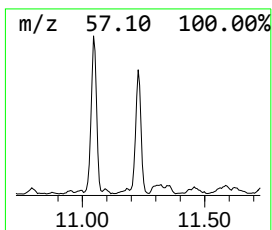
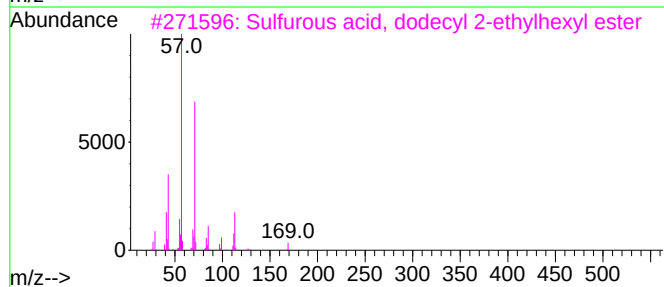
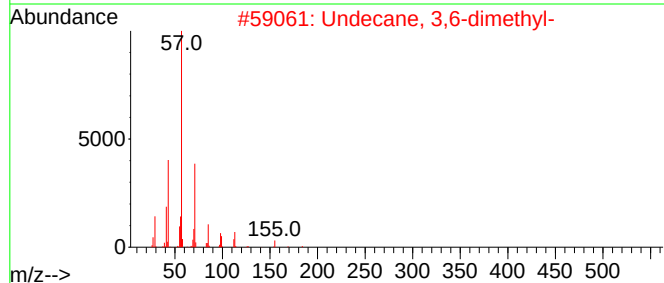
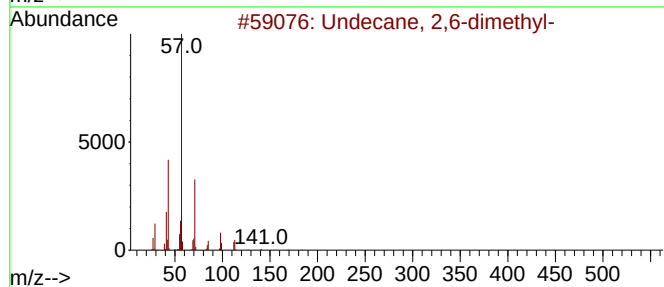
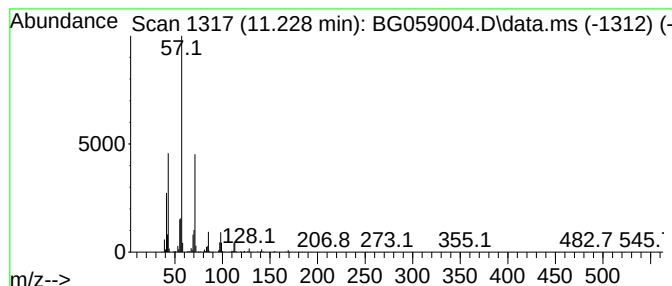
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.228	27.99 ng/ul	2442620	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	70
2			Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	64
3			Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	59
4			Nonane, 4-methyl-	142	C10H22	017301-94-9	59
5			Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

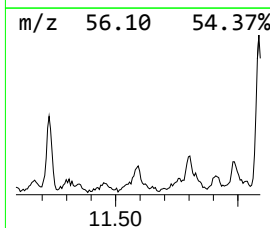
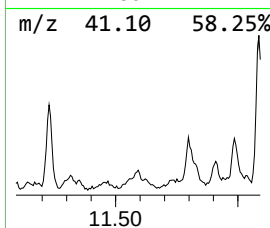
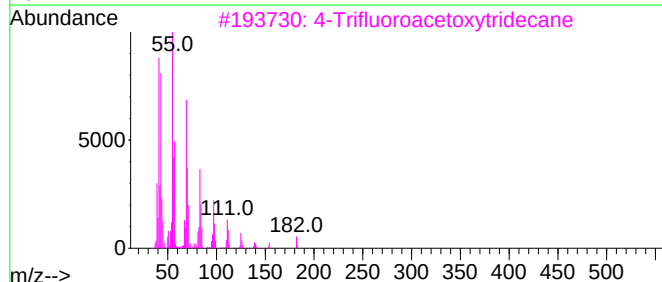
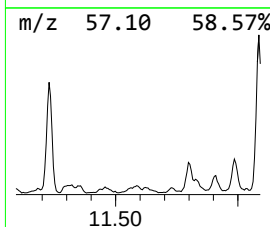
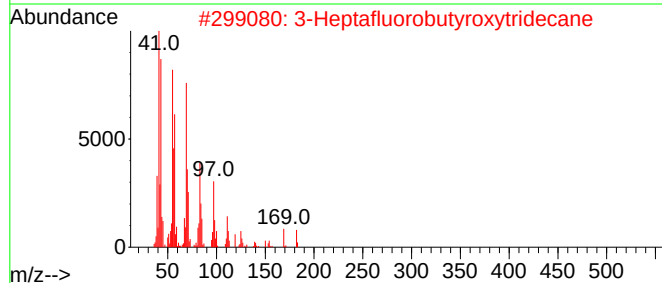
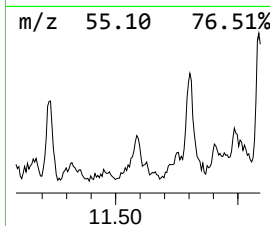
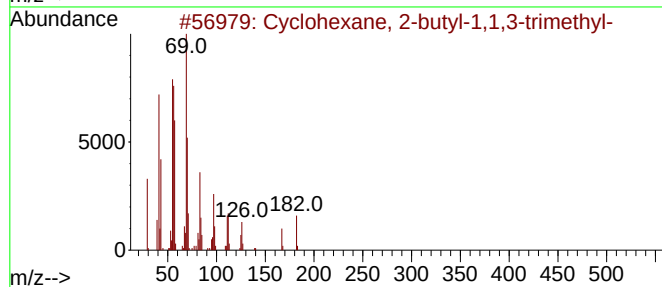
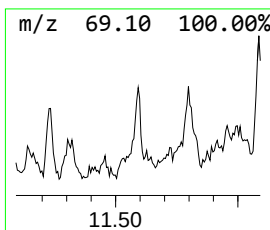
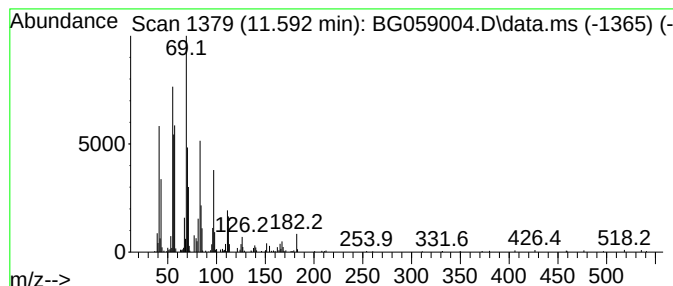
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Cyclic11.592 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.592	11.13 ng/ul	971202	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	90
2	3-Heptafluorobutyroxytridecane	396	C17H27F7O2	1000245-49-5	90
3	4-Trifluoroacetoxytridecane	296	C15H27F3O2	1000245-47-3	90
4	Cyclopentane, 1-pentyl-2-propyl-	182	C13H26	062199-51-3	89
5	3-Tridecene, (E)-	182	C13H26	041446-57-5	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

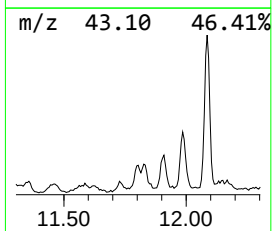
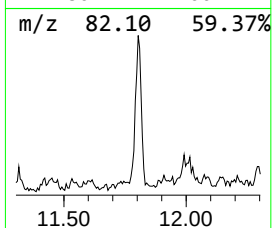
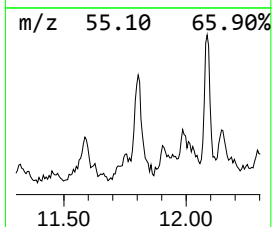
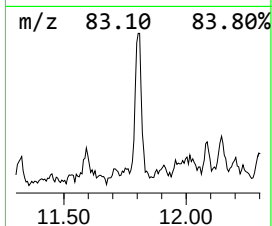
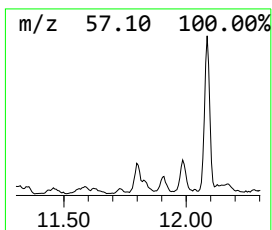
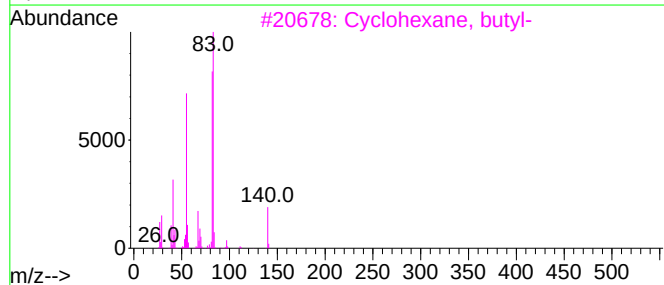
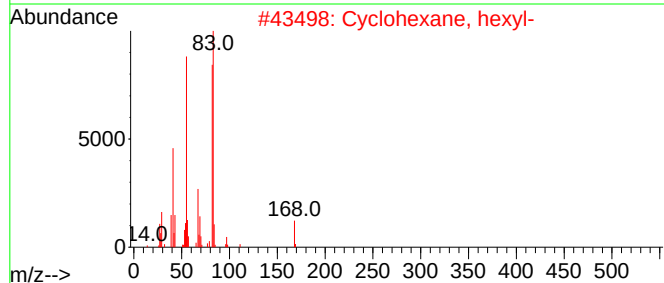
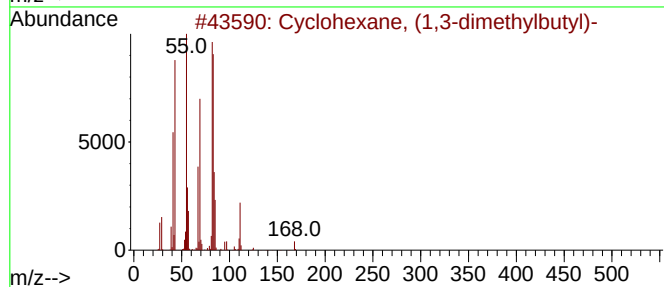
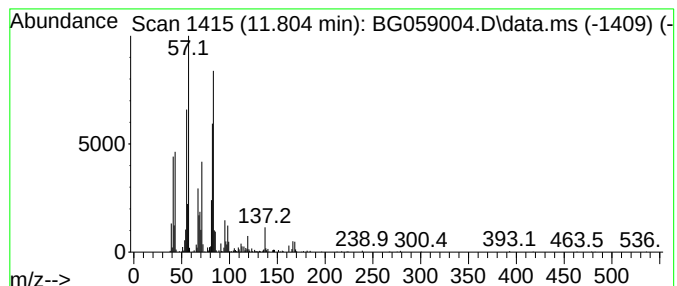
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Cyclic11.804 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.804	23.84 ng/ul	2080480	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1,3-dimethylbutyl)-	168	C12H24	061142-19-6	49
2			Cyclohexane, hexyl-	168	C12H24	004292-75-5	43
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	43
4			Cyclohexane, octyl-	196	C14H28	001795-15-9	43
5			Propanoic acid, 2,2-dimethyl-, c...	184	C11H20O2	029878-49-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

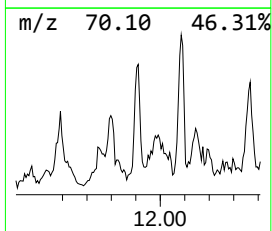
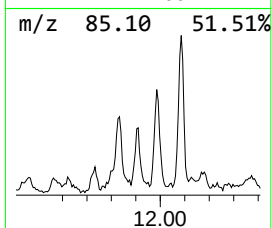
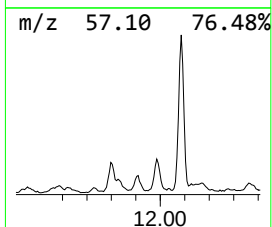
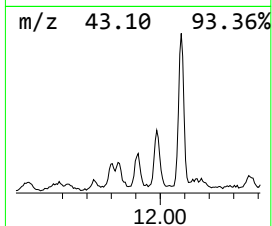
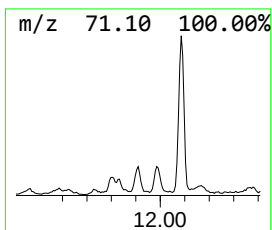
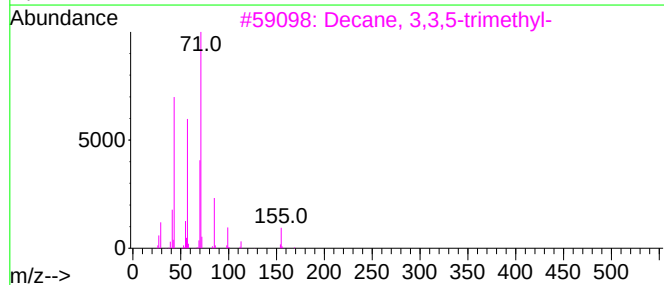
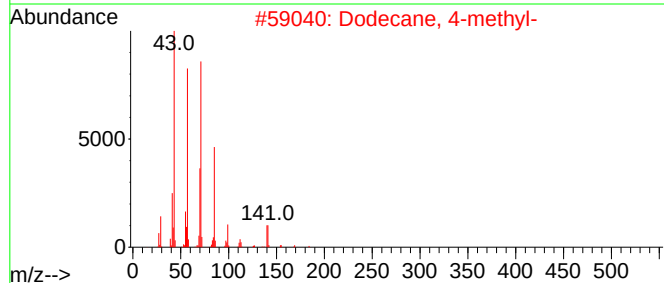
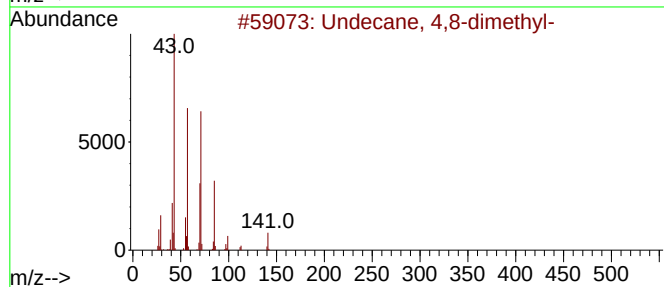
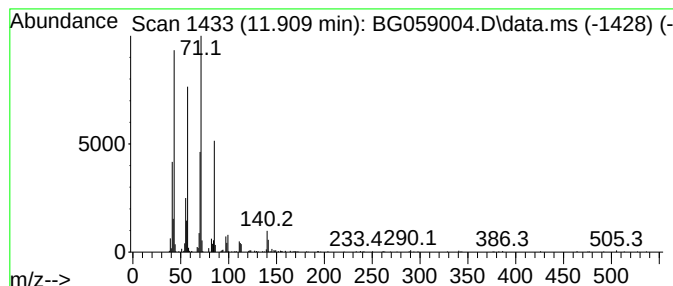
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.909	7.67 ng/ul	669200	Naphthalene-d8	11.169

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	86
2		Dodecane, 4-methyl-	184	C13H28	006117-97-1	86
3		Decane, 3,3,5-trimethyl-	184	C13H28	062338-13-0	64
4		Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	64
5		Nonane, 1-iodo-	254	C9H19I	004282-42-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

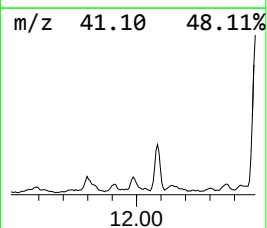
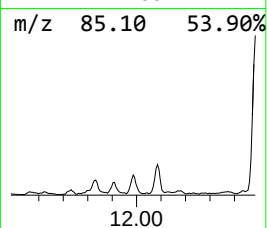
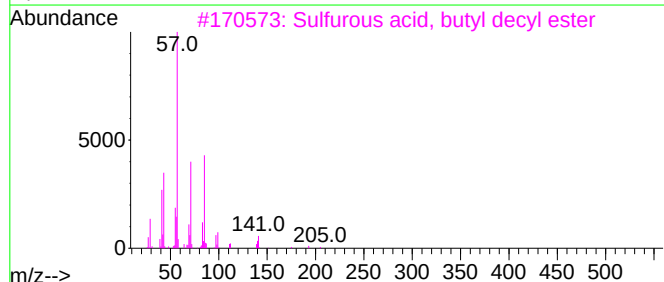
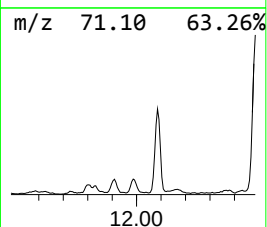
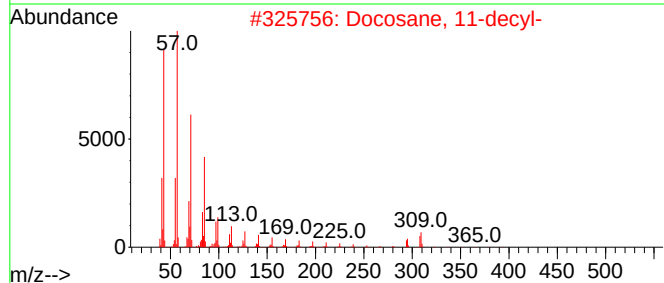
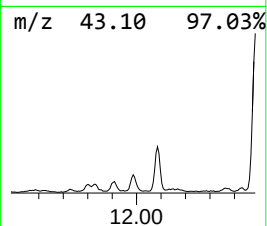
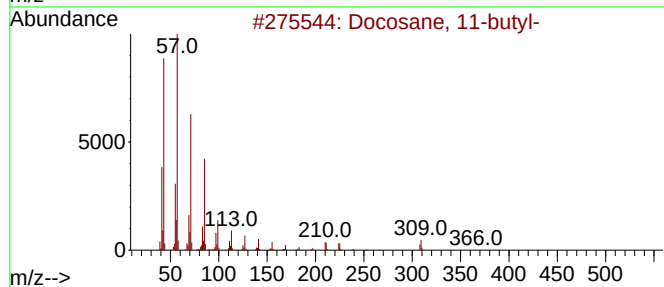
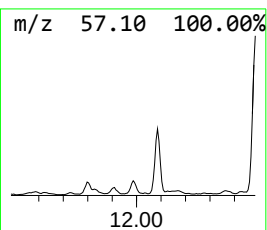
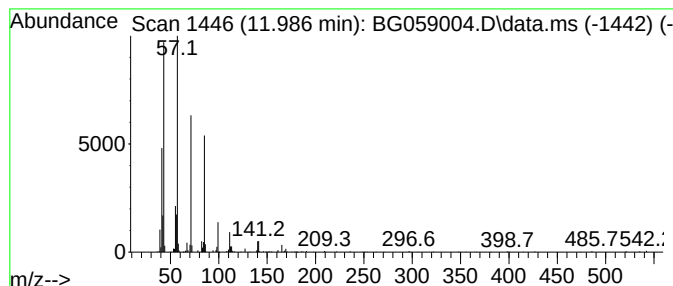
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.986	10.77 ng/ul	939714	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	87
2	Docosane, 11-decyl-	451	C32H66	055401-55-3	78
3	Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	72
4	Undecane, 4-ethyl-	184	C13H28	017312-59-3	72
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

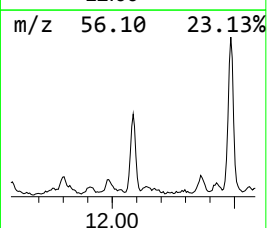
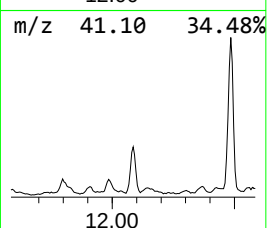
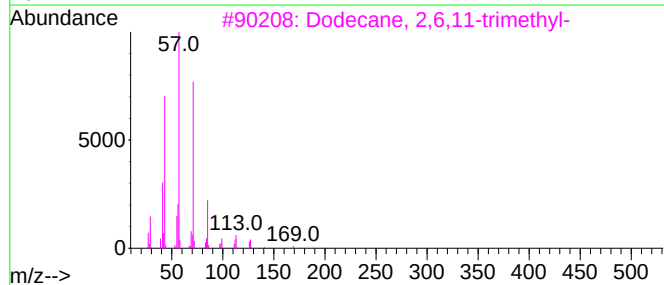
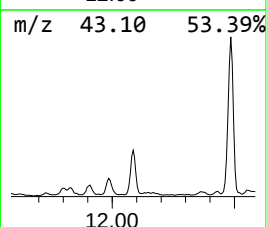
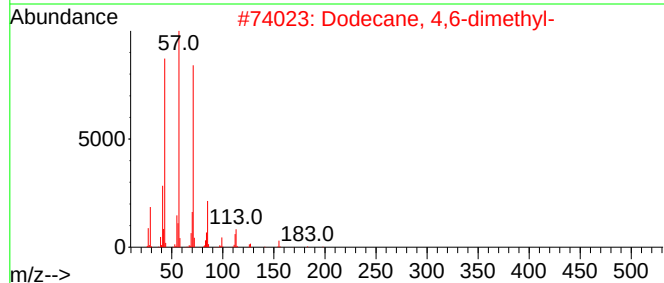
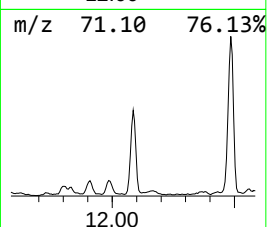
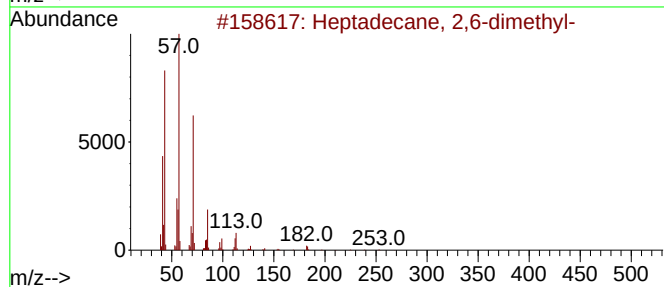
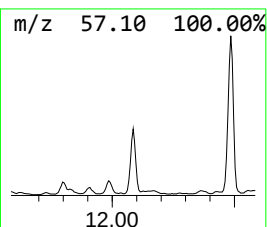
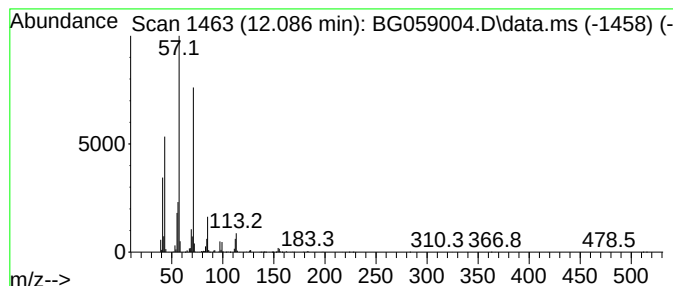
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.086	38.91 ng/ul	3395880	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	83
2	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	83
3	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	78
4	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	72
5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

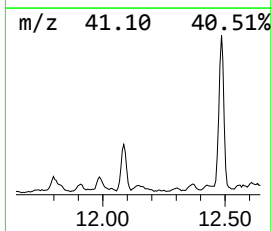
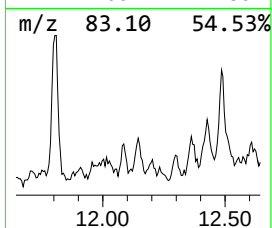
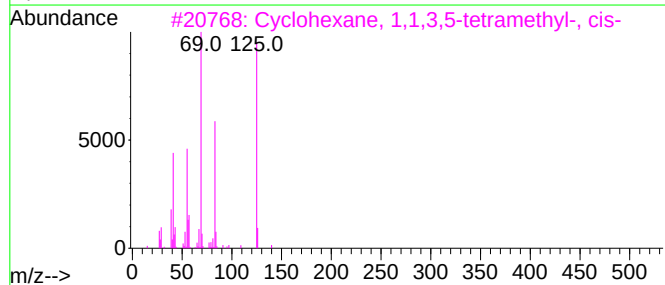
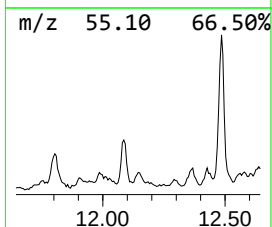
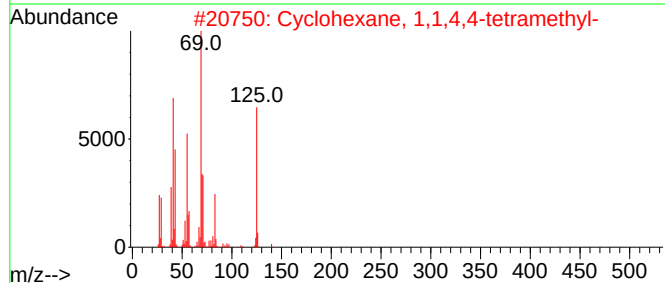
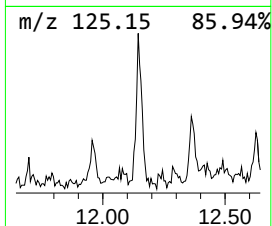
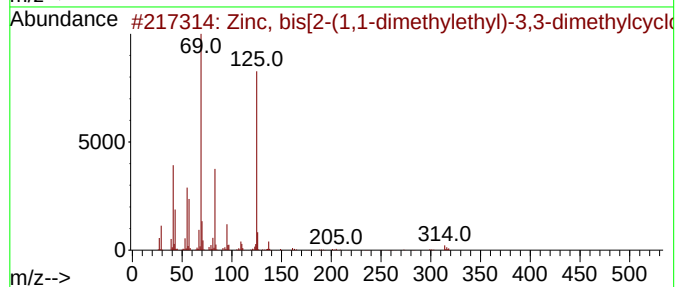
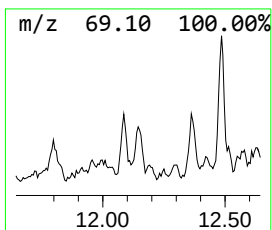
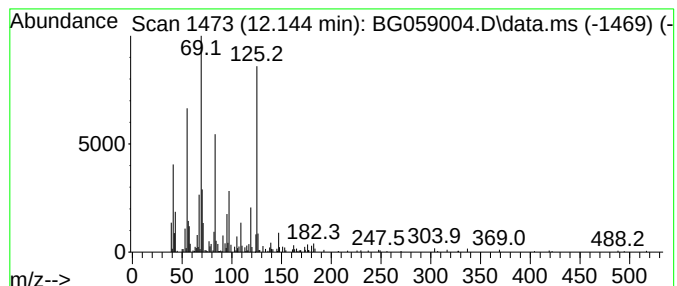
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 Zinc, bis[2-(1,1-dimethylet... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.144	7.24 ng/ul	632199	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Zinc, bis[2-(1,1-dimethylethyl)-...	314	C18H34Zn	074793-36-5	59
2	Cyclohexane, 1,1,4,4-tetramethyl-	140	C10H20	002223-52-1	58
3	Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-32-9	58
4	Cyclohexane, 1,1,3,5-tetramethyl...	140	C10H20	050876-31-8	50
5	Cyclopentane, 1,2-dibutyl-	182	C13H26	062199-52-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

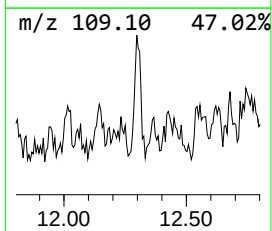
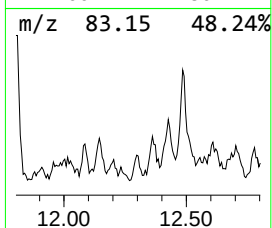
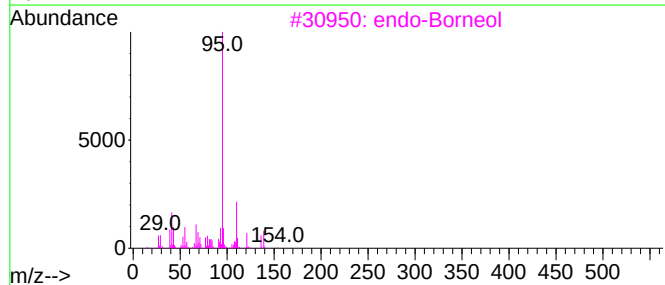
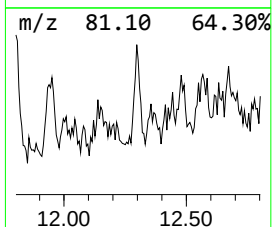
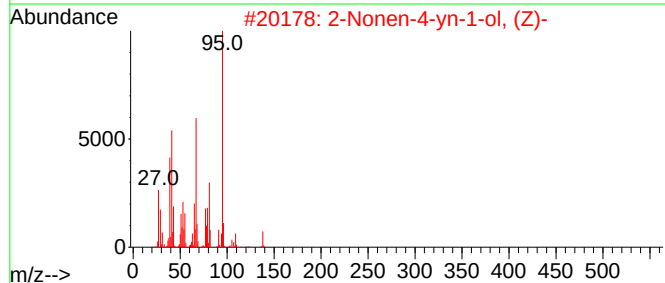
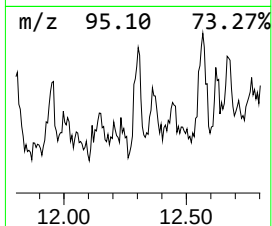
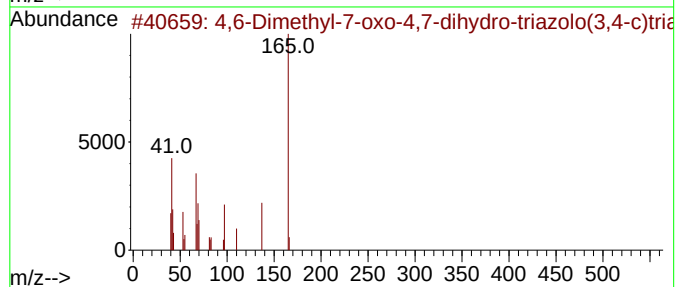
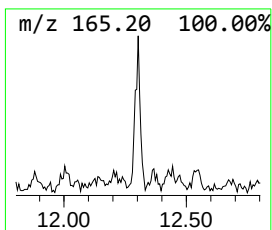
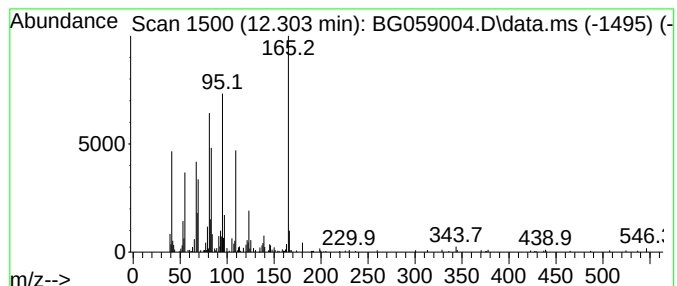
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 unknown-04

Concentration Rank 72

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.303	4.36 ng/ul	380868	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,6-Dimethyl-7-oxo-4,7-dihydro-t...	165	C6H7N5O	025623-69-2	27
2			2-Nonen-4-yn-1-ol, (Z)-	138	C9H14O	134225-90-4	27
3			endo-Borneol	154	C10H18O	000507-70-0	22
4			exo-2-Bromonorbornane	174	C7H11Br	002534-77-2	22
5			2-Norbornyl bromide	174	C7H11Br	029342-65-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

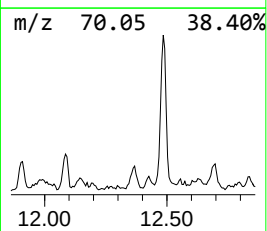
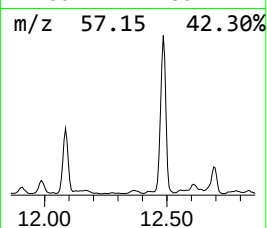
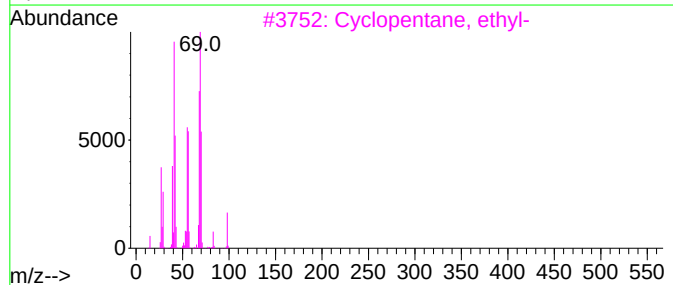
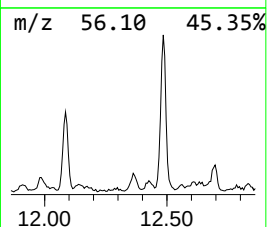
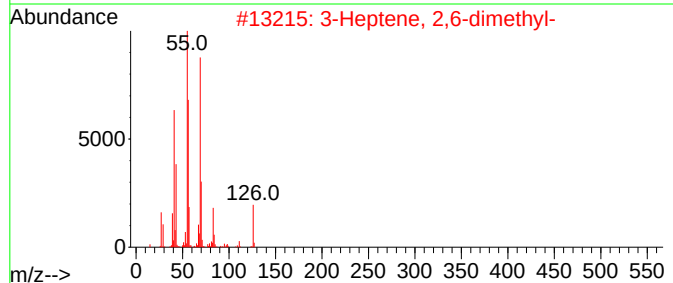
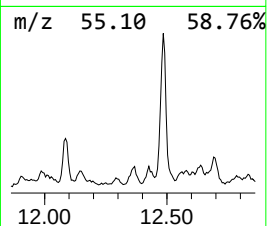
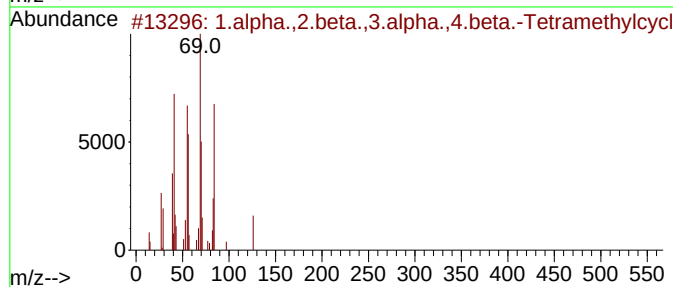
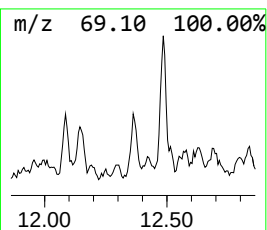
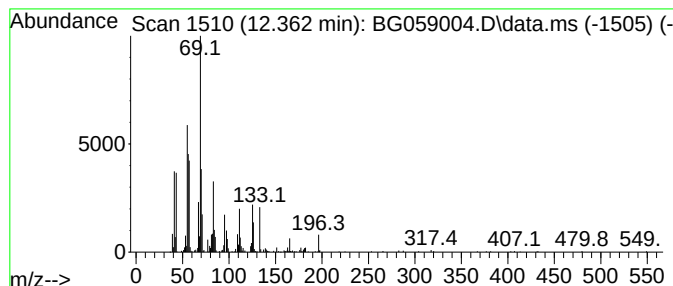
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.362	9.17 ng/ul	800430	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1.alpha.,2.beta.,3.alpha.,4.beta...			126	C9H18	002532-67-4	58
2	3-Heptene, 2,6-dimethyl-			126	C9H18	002738-18-3	53
3	Cyclopentane, ethyl-			98	C7H14	001640-89-7	52
4	1-Hexene, 3,3-dimethyl-			112	C8H16	003404-77-1	49
5	trans-2-Methyl-3-octene			126	C9H18	052937-36-7	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

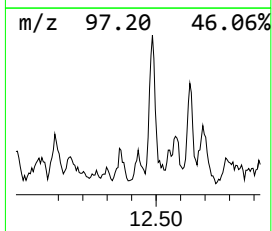
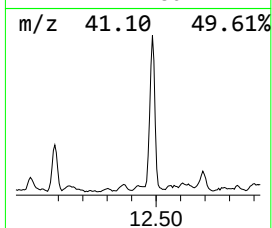
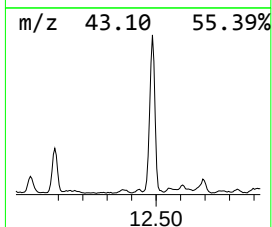
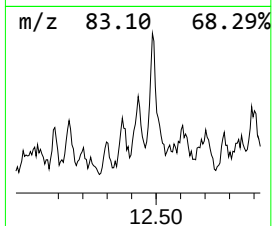
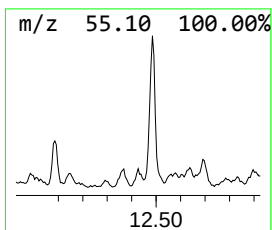
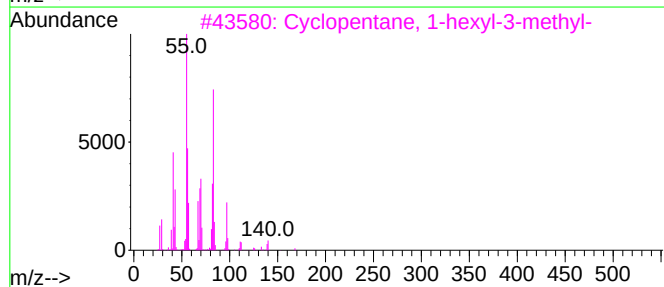
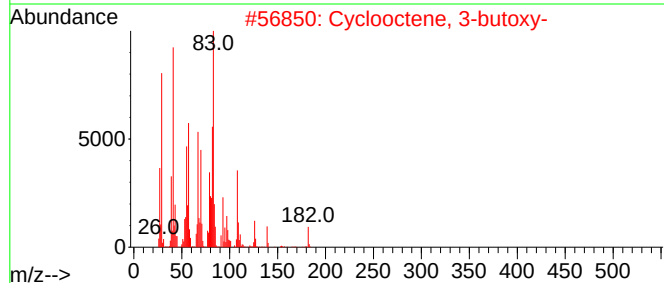
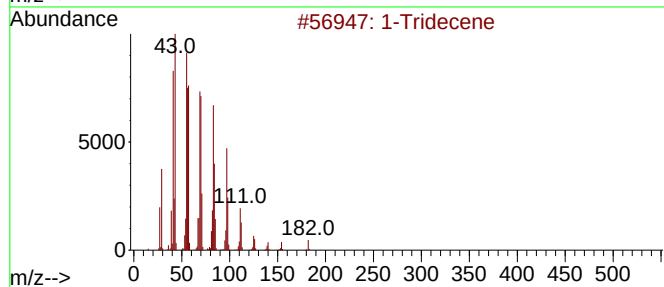
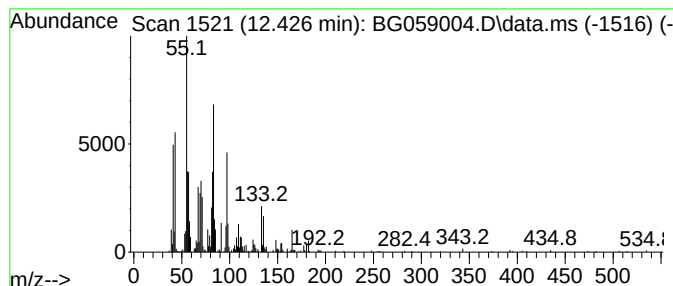
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.426	8.78 ng/ul	766116	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tridecene	182	C13H26	002437-56-1	46
2	Cyclooctene, 3-butoxy-	182	C12H22O	1000159-08-1	30
3	Cyclopentane, 1-hexyl-3-methyl-	168	C12H24	061142-68-5	25
4	1-Tridecene	182	C13H26	002437-56-1	25
5	Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

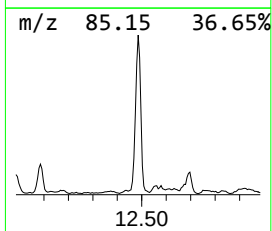
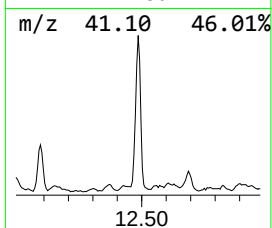
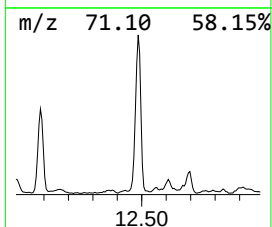
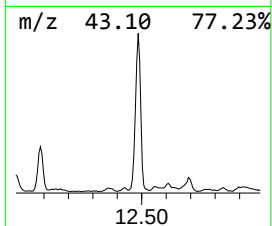
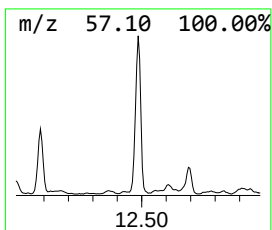
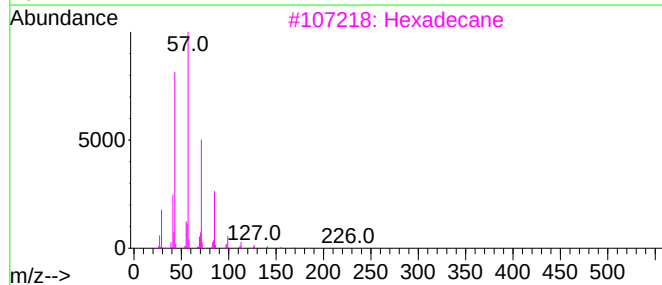
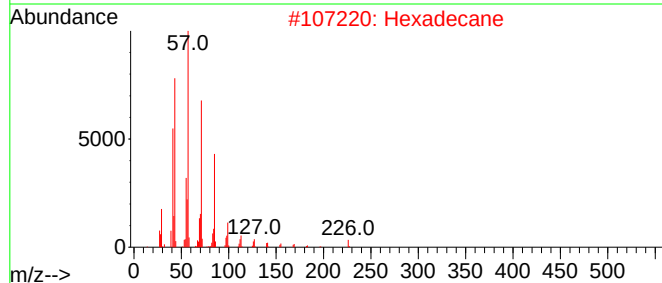
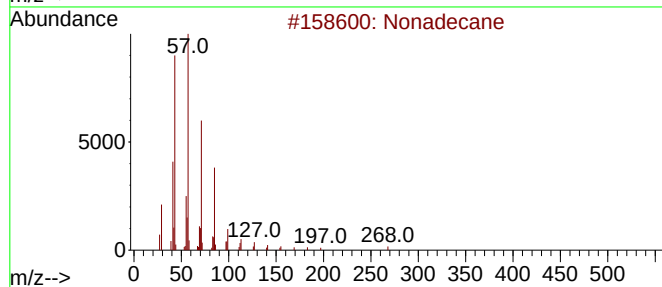
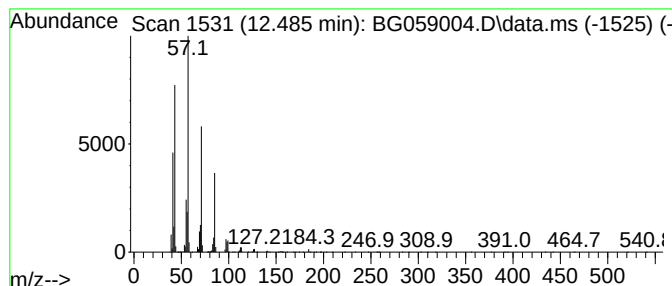
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.485	110.96 ng/ul	9684610	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	86
2	Hexadecane	226	C16H34	000544-76-3	86
3	Hexadecane	226	C16H34	000544-76-3	86
4	Tridecane	184	C13H28	000629-50-5	86
5	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

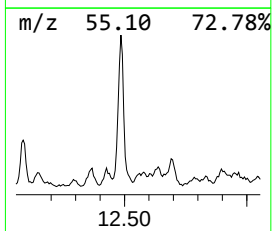
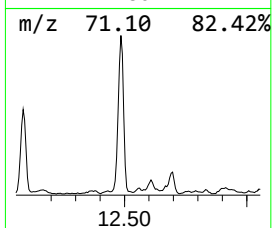
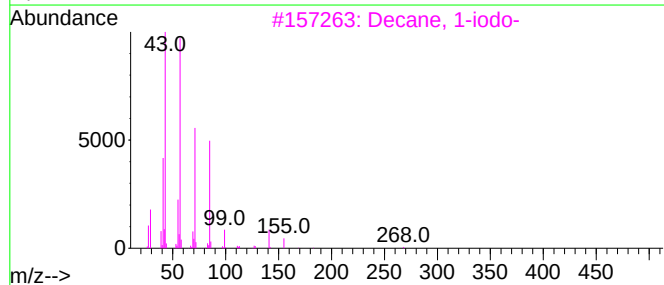
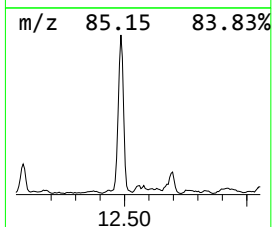
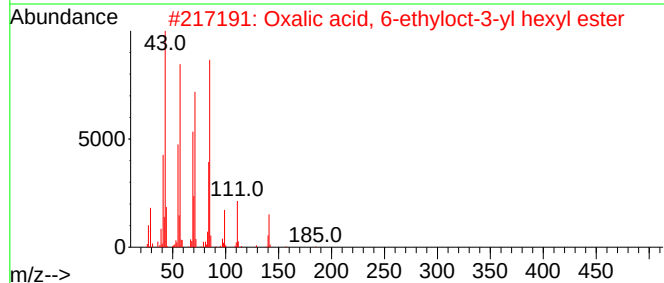
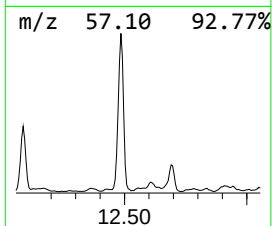
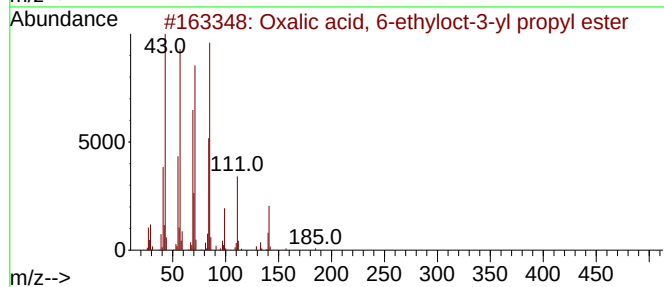
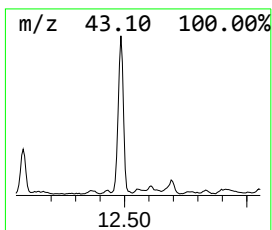
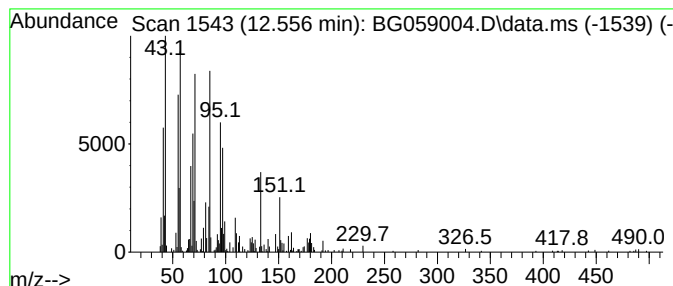
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 unknown-05 Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.556	7.58 ng/ul	661293	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 6-ethyloct-3-yl pro...	272	C15H28O4	1000309-34-0	38
2			Oxalic acid, 6-ethyloct-3-yl hex...	314	C18H34O4	1000309-34-4	38
3			Decane, 1-iodo-	268	C10H21I	002050-77-3	35
4			Oxalic acid, 6-ethyloct-3-yl iso...	314	C18H34O4	1000309-34-3	35
5			Oxalic acid, butyl 6-ethyloct-3-...	286	C16H30O4	1000309-34-2	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

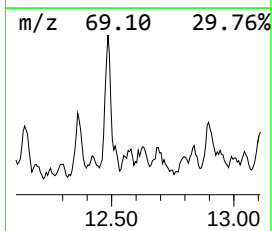
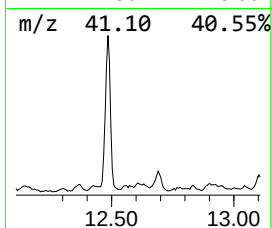
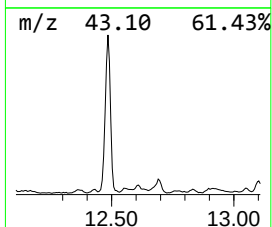
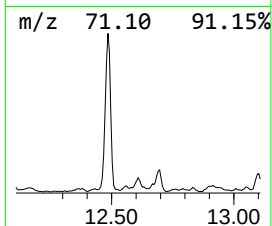
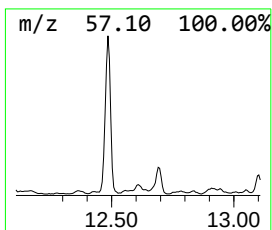
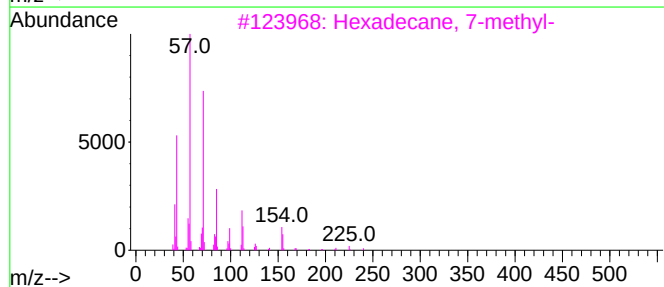
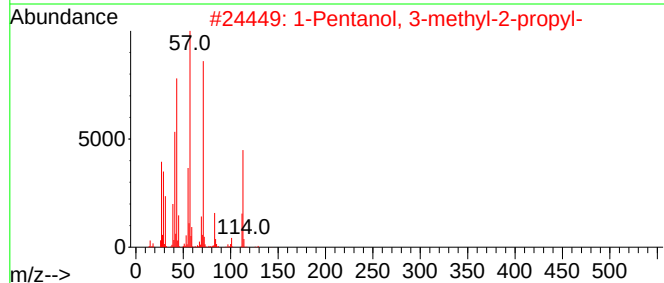
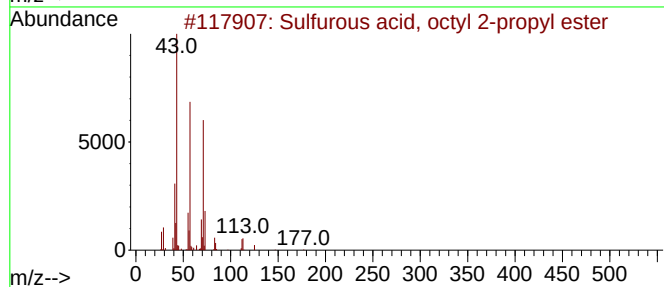
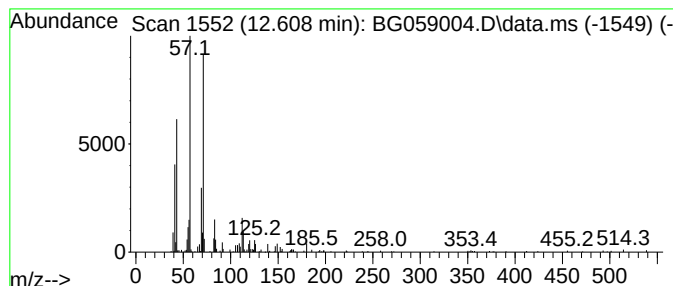
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 Sulfurous acid, octyl 2-pro... Concentration Rank 76

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.608	3.60 ng/ul	314056	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, octyl 2-propyl e...	236	C11H24O3S	1000309-11-9	78
2			1-Pentanol, 3-methyl-2-propyl-	144	C9H20O	054004-40-9	72
3			Hexadecane, 7-methyl-	240	C17H36	026730-20-1	72
4			Nonane, 5-propyl-	170	C12H26	000998-35-6	72
5			Octane, 2-bromo-	192	C8H17Br	000557-35-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

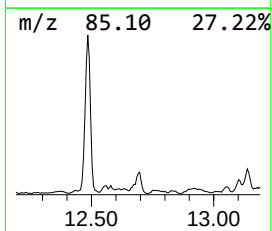
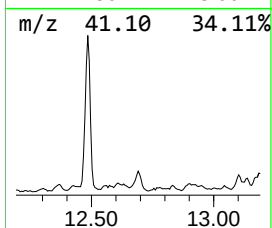
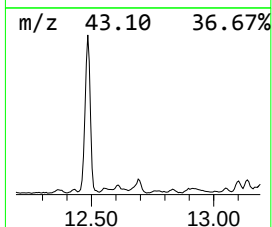
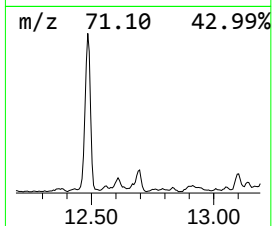
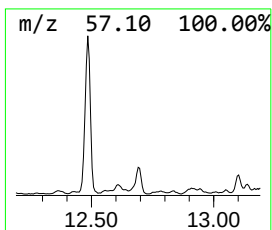
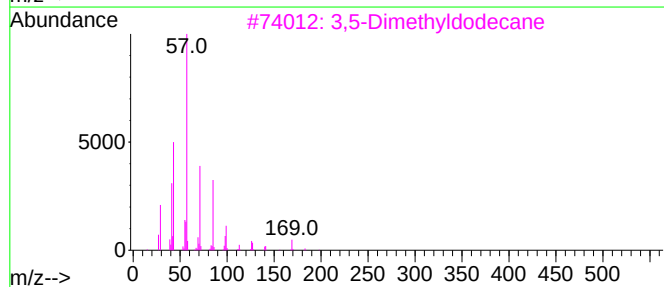
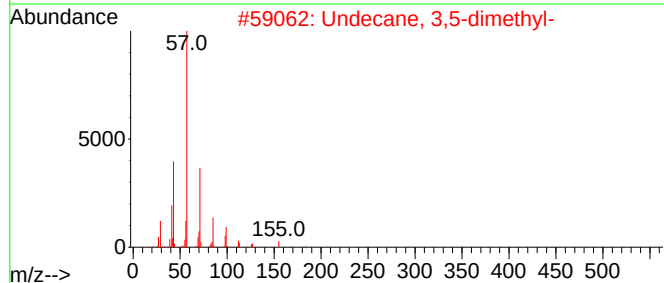
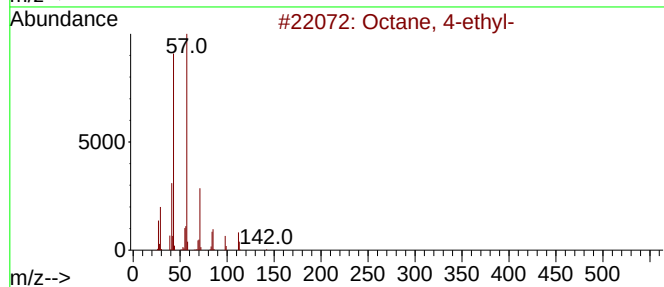
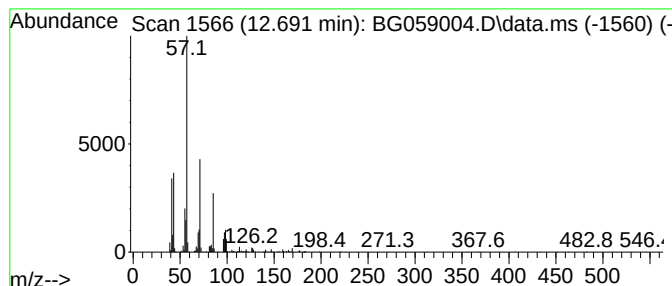
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.691	21.79 ng/ul	1901780	Naphthalene-d8	11.169	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-ethyl-	142	C10H22	015869-86-0	64
2	Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	59
3	3,5-Dimethyldodecane	198	C14H30	107770-99-0	53
4	Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	53
5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

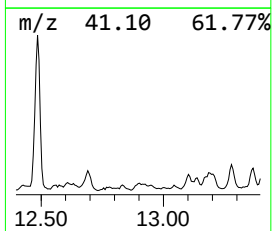
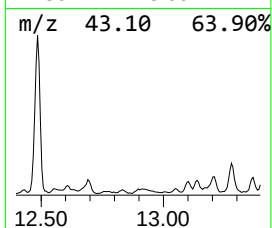
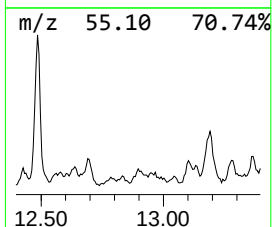
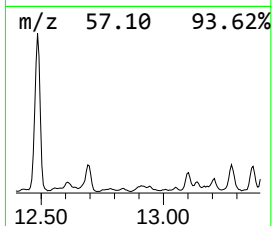
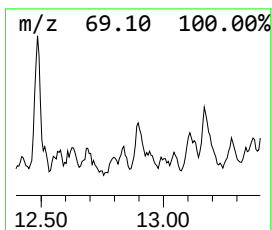
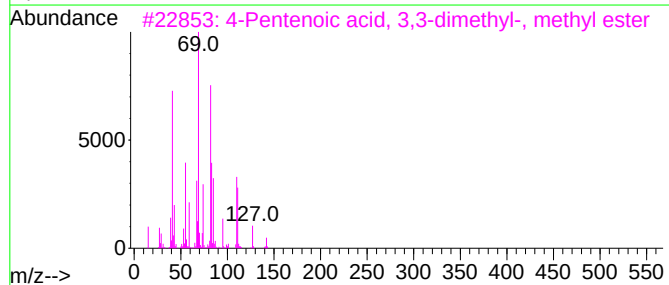
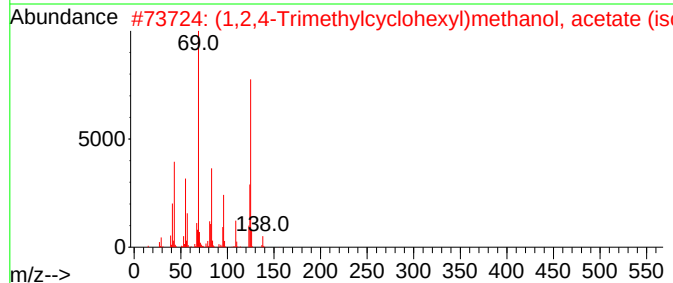
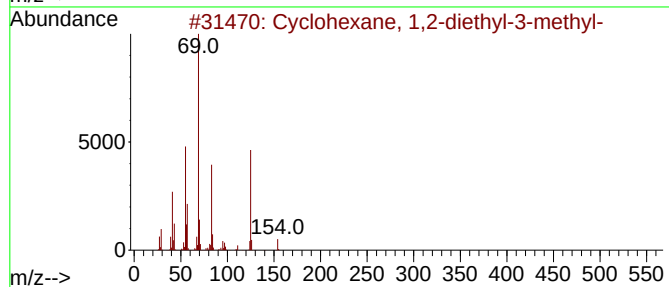
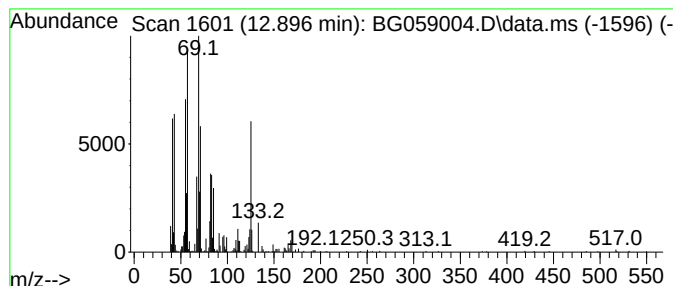
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 (DEL) Alkane: Cyclic12.896 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.896	12.54 ng/ul	1094700	Naphthalene-d8	11.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-diethyl-3-methyl-	154	C11H22	061141-80-8	47
2			(1,2,4-Trimethylcyclohexyl)metha...	198	C12H22O2	1000499-04-5	47
3			4-Pentenoic acid, 3,3-dimethyl-,...	142	C8H14O2	063721-05-1	27
4			3-Methyl-2-hexene	98	C7H14	017618-77-8	27
5			1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

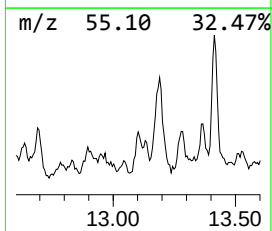
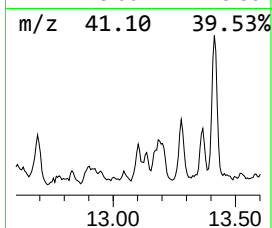
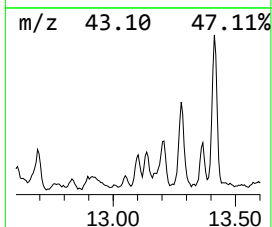
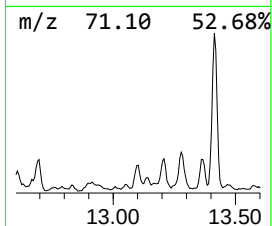
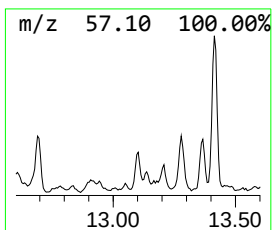
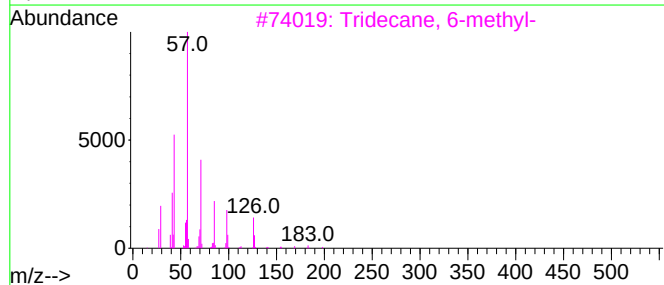
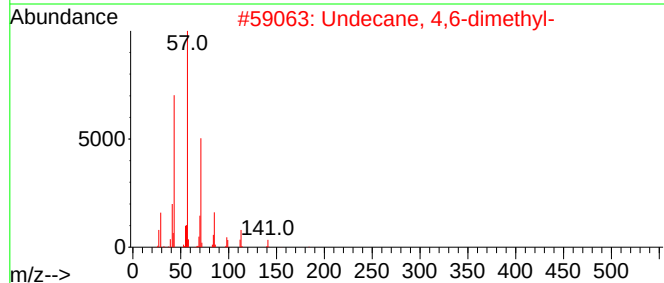
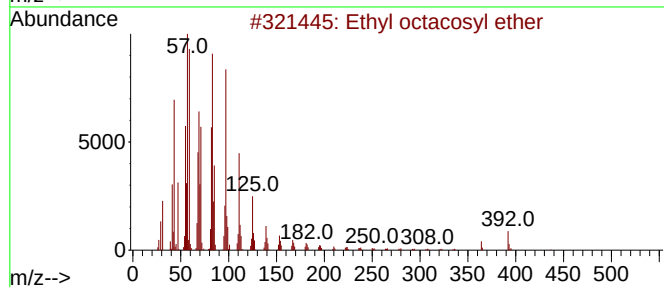
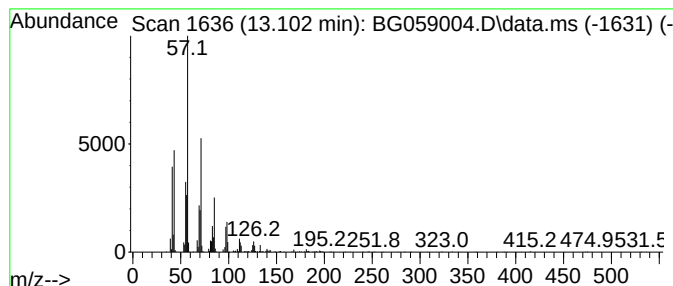
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 Ethyl octacosyl ether Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.102	22.97 ng/ul	1218560	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl octacosyl ether	438	C30H62O	1010406-37-0	80
2	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	58
3	Tridecane, 6-methyl-	198	C14H30	013287-21-3	53
4	Hexadecane	226	C16H34	000544-76-3	50
5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

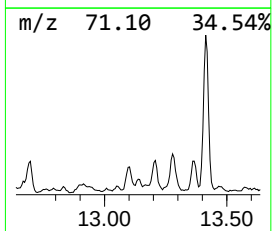
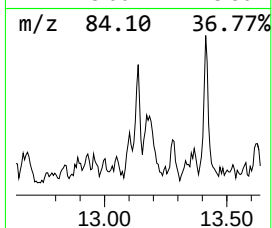
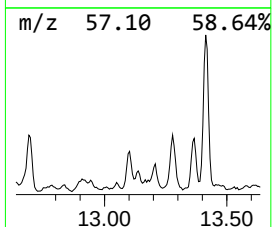
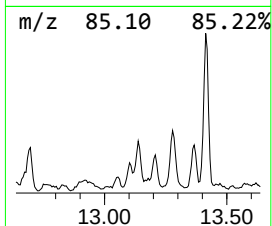
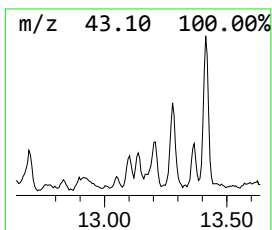
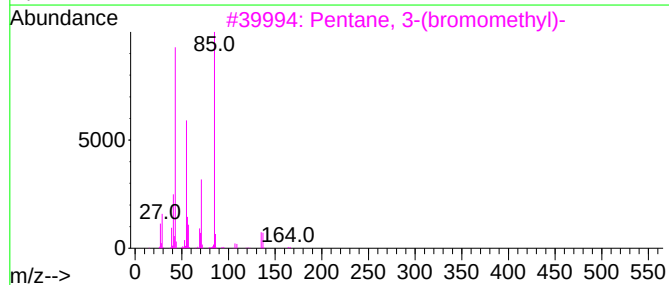
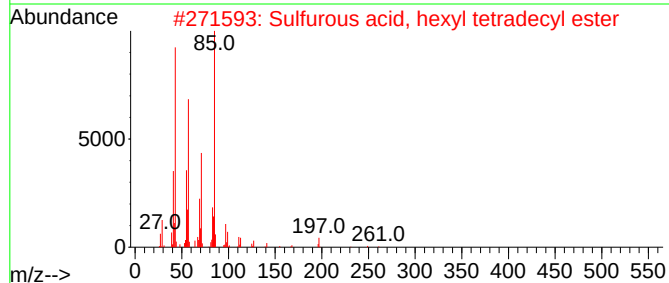
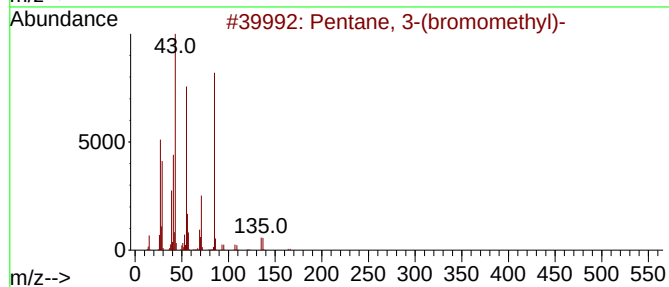
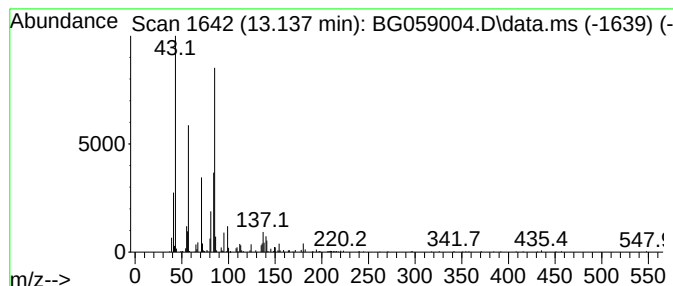
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 Pentane, 3-(bromomethyl)- Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.137	15.48 ng/ul	821173	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 3-(bromomethyl)-	164	C6H13Br	003814-34-4	53
2			Sulfurous acid, hexyl tetradecyl...	362	C20H42O3S	1010309-13-6	53
3			Pentane, 3-(bromomethyl)-	164	C6H13Br	003814-34-4	53
4			Silane, ethenylethylidimethyl-	114	C6H14Si	018163-06-9	50
5			Pentane, 2,3,3,4-tetramethyl-	128	C9H20	016747-38-9	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

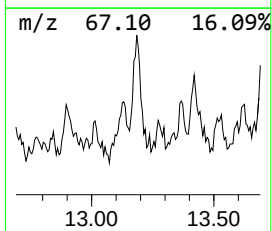
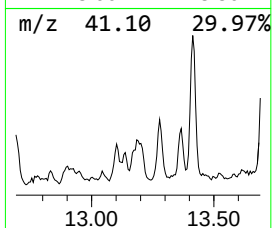
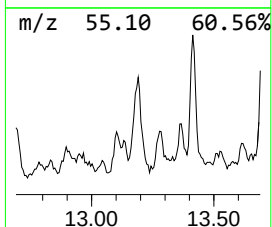
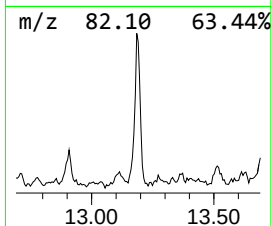
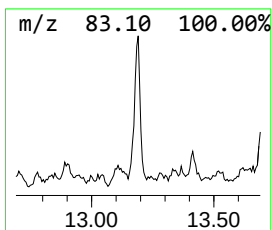
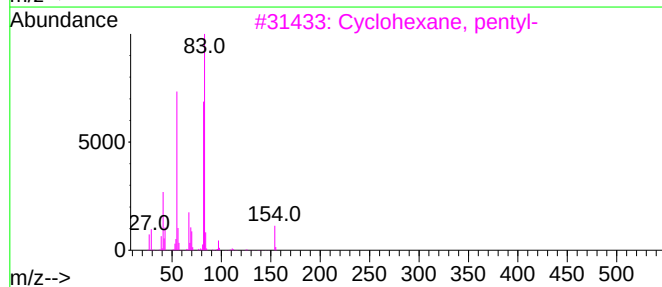
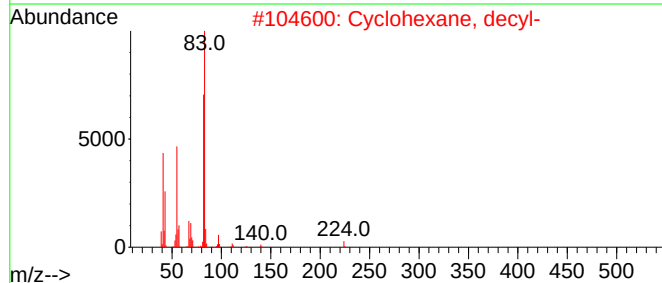
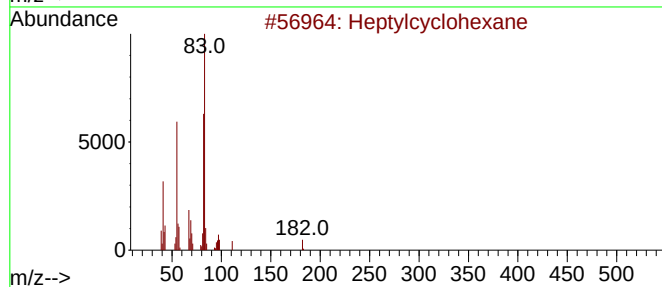
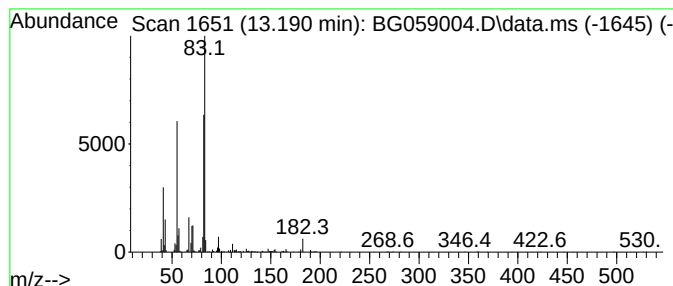
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 (DEL) Alkane: Cyclic13.190 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.190	48.46 ng/ul	2571370	Acenaphthene-d10		14.947	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptylcyclohexane		182	C13H26	005617-41-4	80
2	Cyclohexane, decyl-		224	C16H32	001795-16-0	72
3	Cyclohexane, pentyl-		154	C11H22	004292-92-6	64
4	Cyclohexane, pentyl-		154	C11H22	004292-92-6	64
5	Cyclohexane, octyl-		196	C14H28	001795-15-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

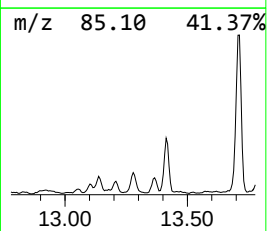
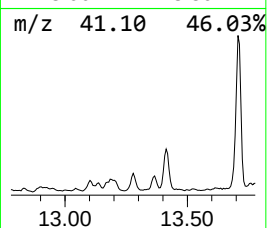
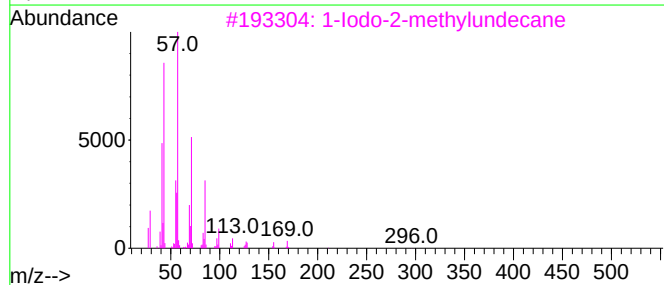
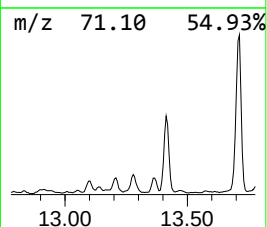
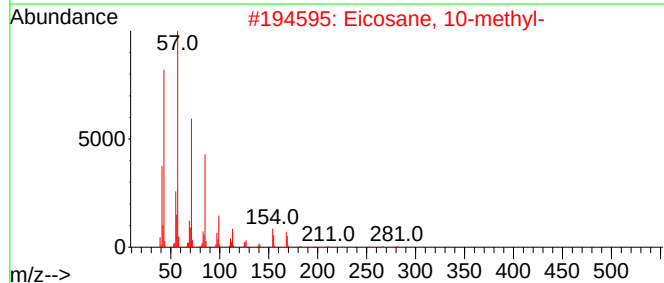
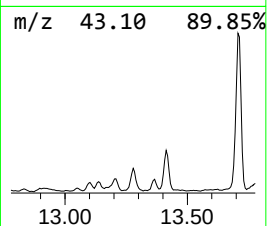
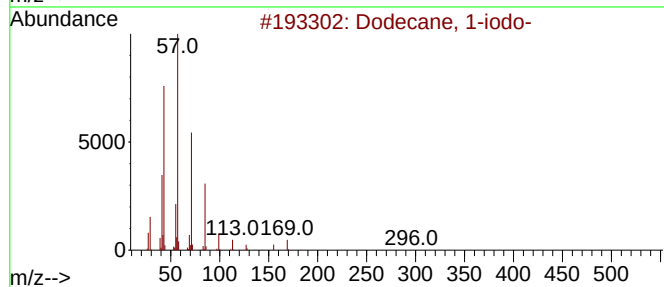
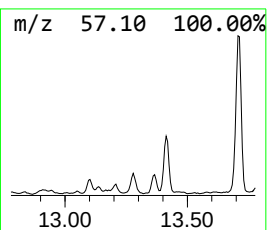
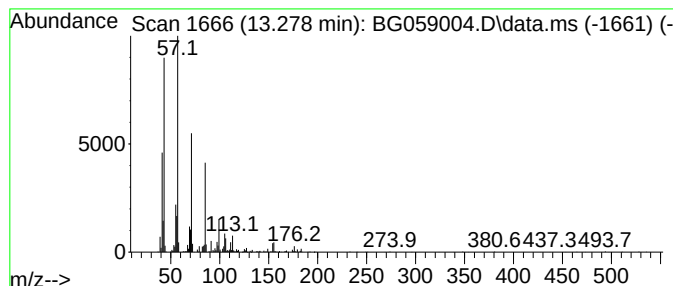
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 Dodecane, 1-iodo- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.278	34.94 ng/ul	1854170	Acenaphthene-d10	14.947

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	94
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	87
4		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5		Nonadecane, 2-methyl-	282	C20H42	001560-86-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

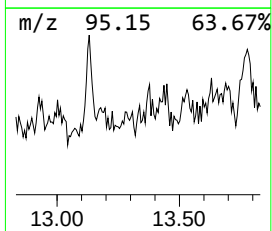
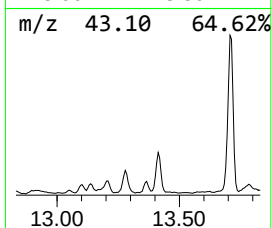
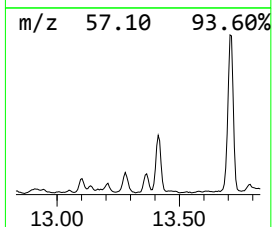
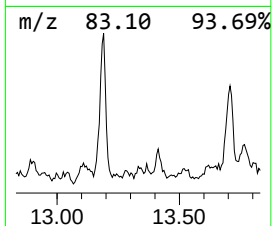
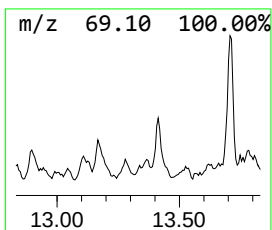
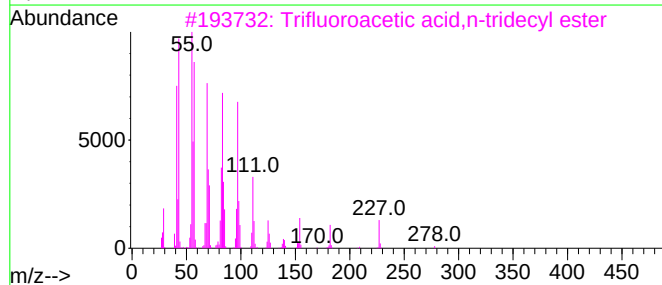
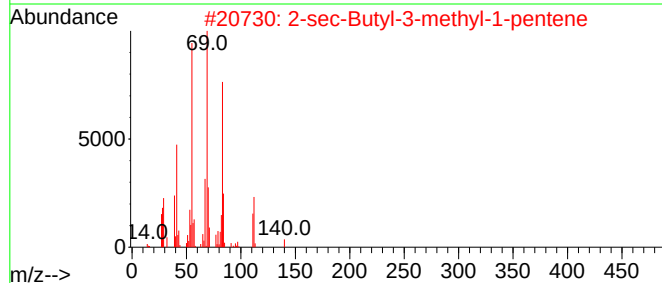
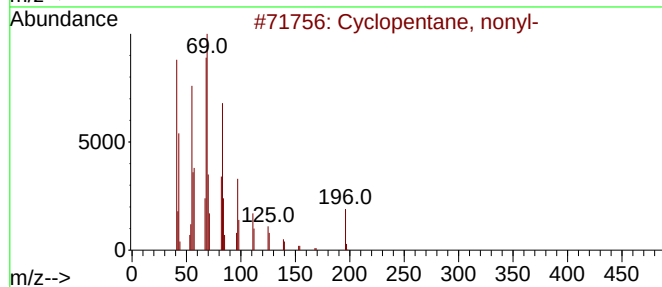
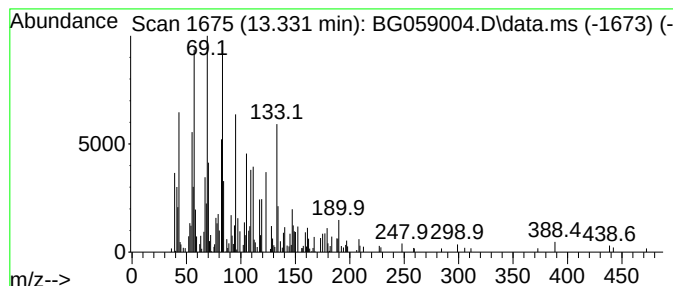
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 (DEL) Alkane: Cyclic13.331 Concentration Rank 80

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.331	2.66 ng/ul	141071	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, nonyl-	196	C14H28	002882-98-6	46
2			2-sec-Butyl-3-methyl-1-pentene	140	C10H20	075144-24-0	45
3			Trifluoroacetic acid,n-tridecyl ...	296	C15H27F3O2	053800-02-5	38
4			n-Heptadecanol-1	256	C17H36O	001454-85-9	38
5			Bicyclo[3.1.1]heptan-3-one, 2,6,...	152	C10H16O	015358-88-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

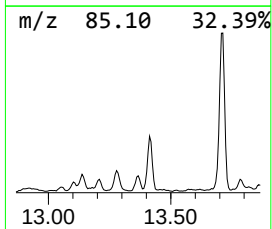
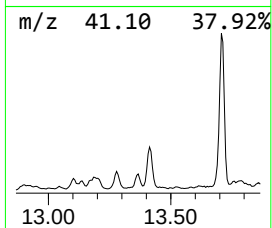
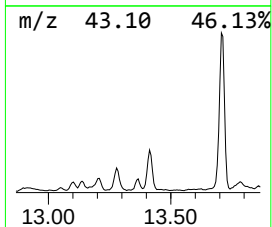
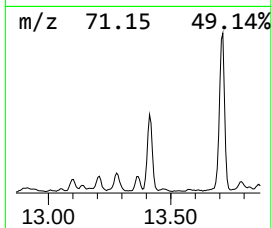
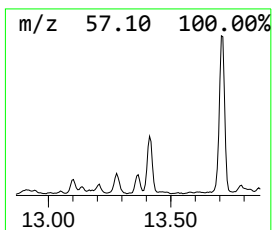
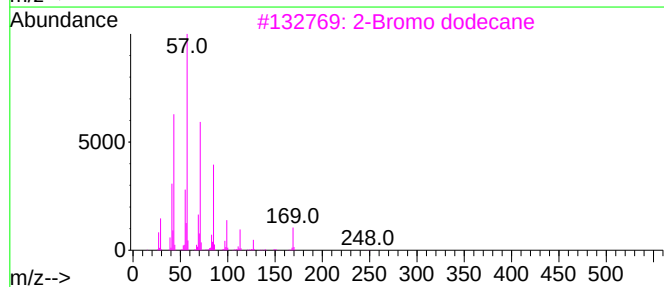
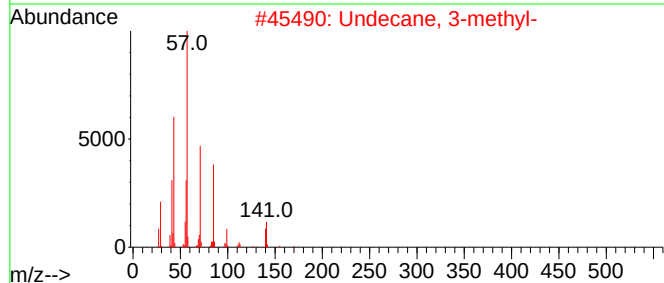
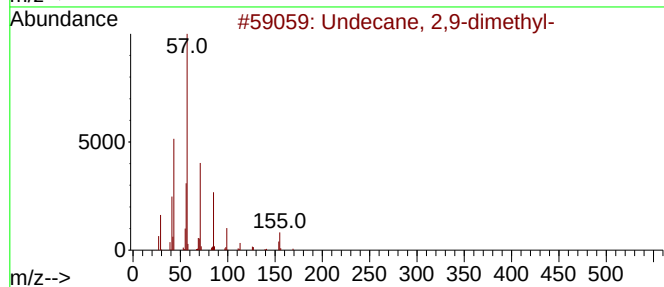
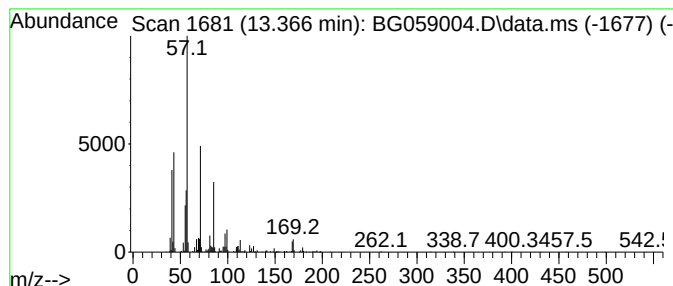
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.366	24.10 ng/ul	1278530	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	80
2	Undecane, 3-methyl-	170	C12H26	001002-43-3	80
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	74
4	Undecane, 3-methyl-	170	C12H26	001002-43-3	72
5	Tridecane, 2-methyl-	198	C14H30	001560-96-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

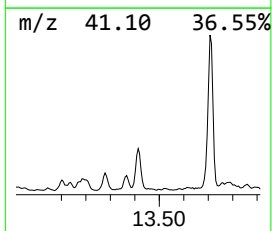
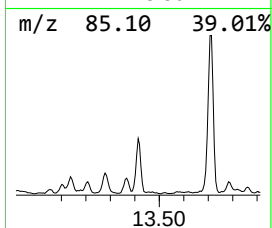
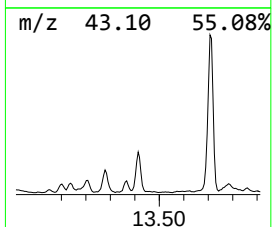
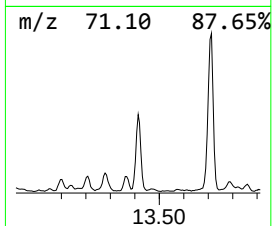
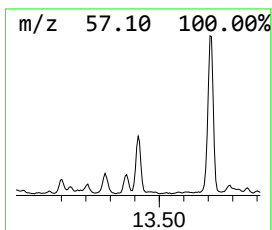
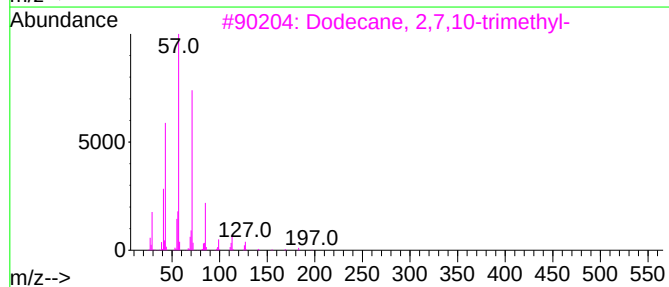
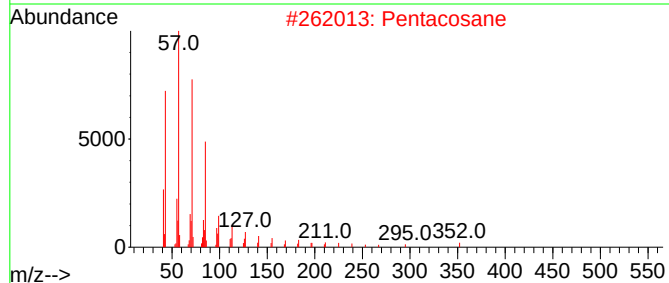
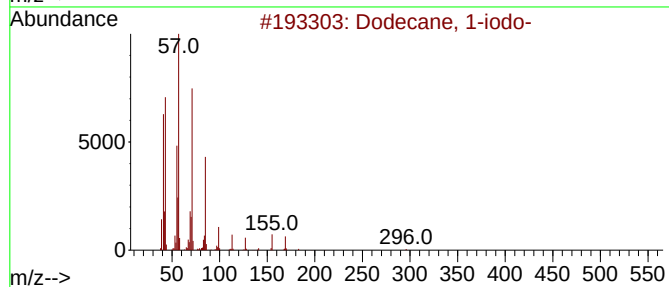
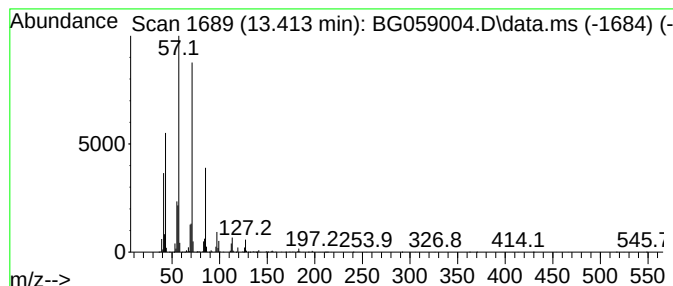
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 Pentacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.413	83.37 ng/ul	4423480	Acenaphthene-d10		14.947	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	86
2	Pentacosane		352	C25H52	000629-99-2	86
3	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	80
4	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	80
5	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

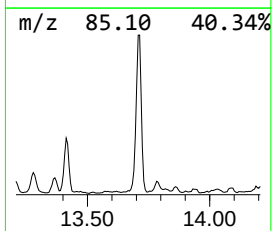
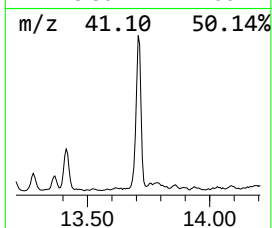
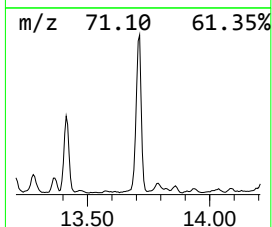
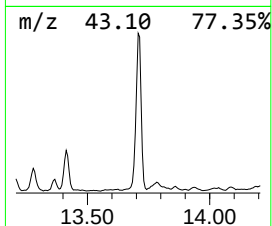
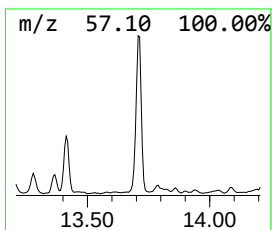
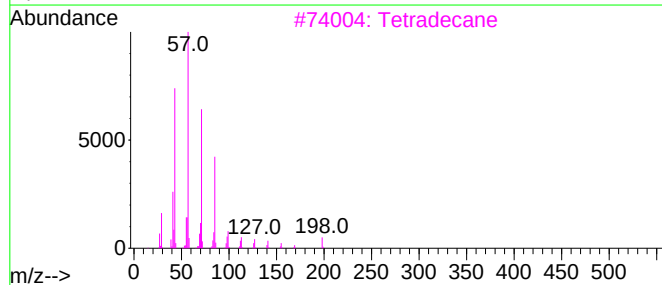
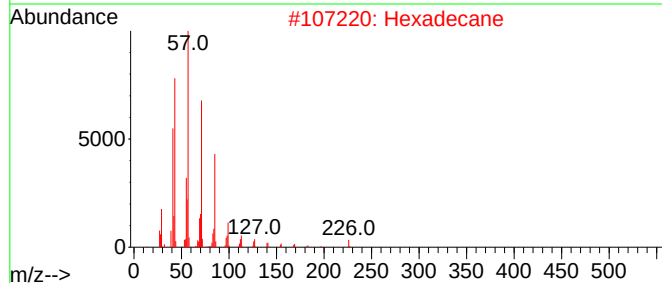
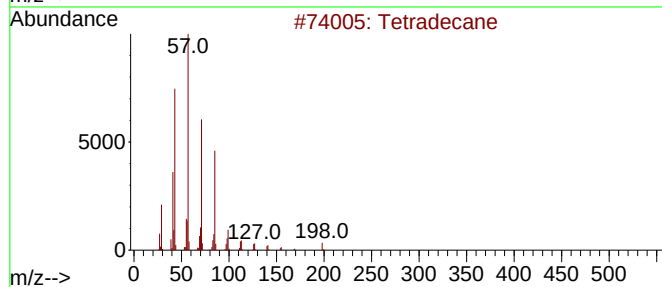
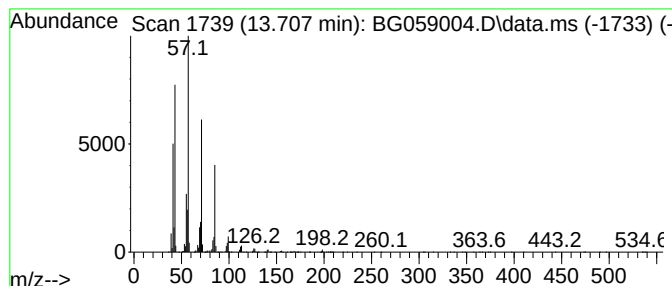
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.707	252.46 ng/ul	13395800	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	95
2	Hexadecane	226	C16H34	000544-76-3	90
3	Tetradecane	198	C14H30	000629-59-4	87
4	Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	86
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

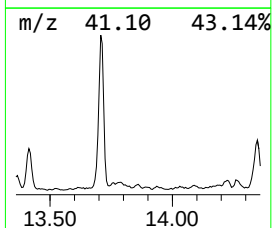
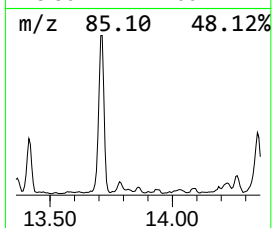
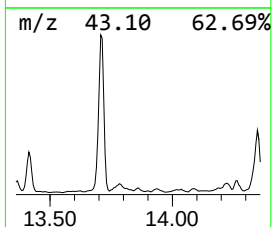
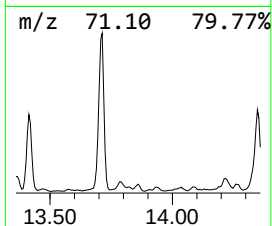
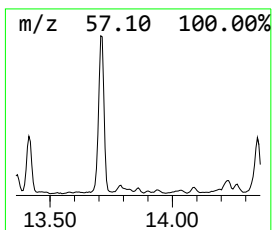
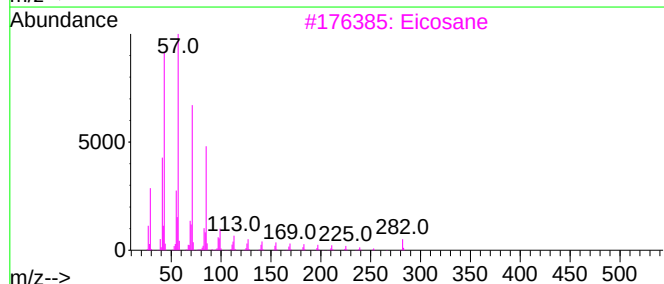
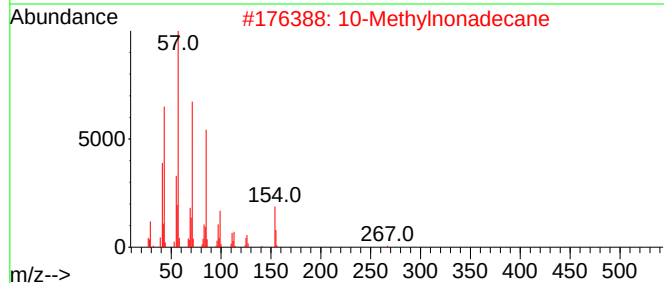
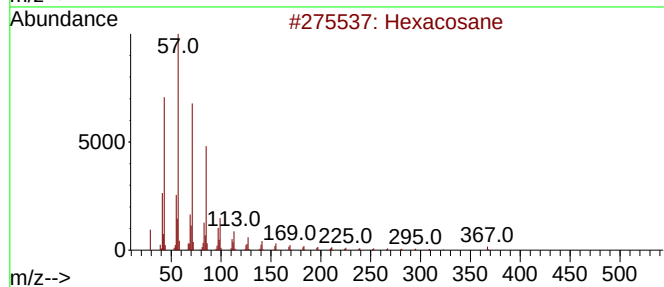
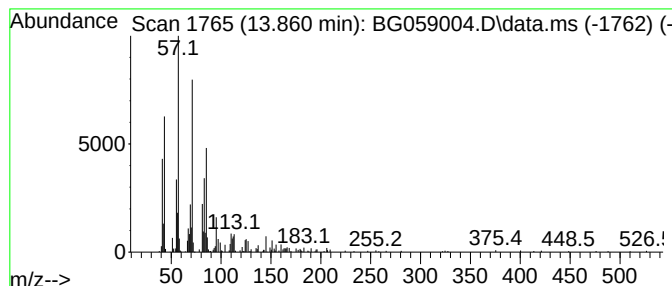
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.860	6.46 ng/ul	343008	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	72
2			10-Methylnonadecane	282	C20H42	056862-62-5	72
3			Eicosane	282	C20H42	000112-95-8	72
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72
5			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

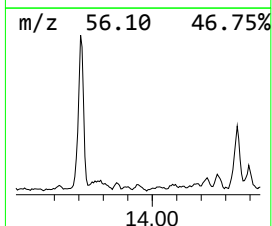
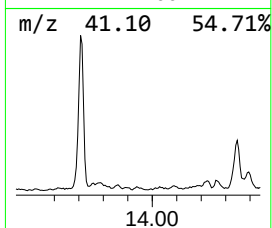
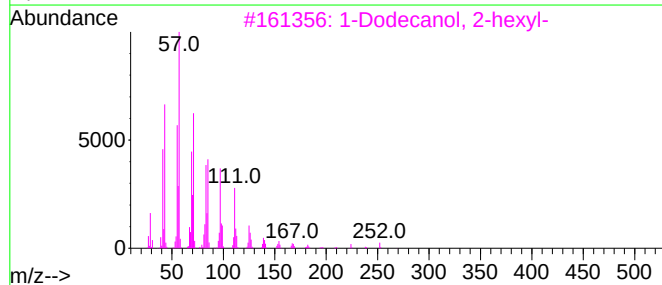
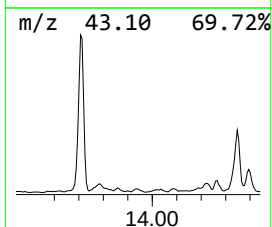
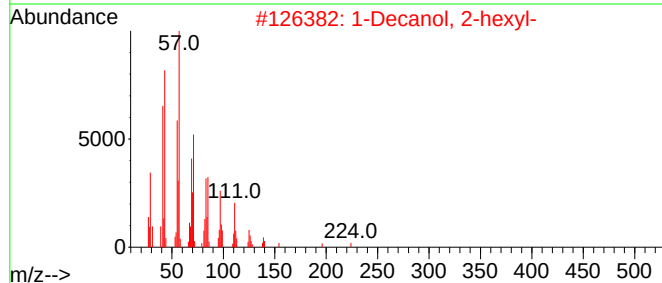
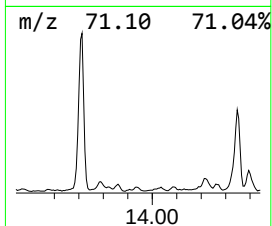
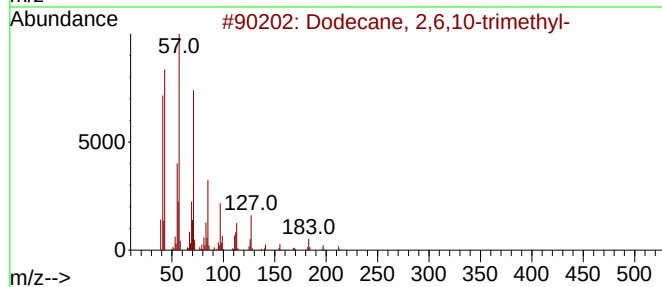
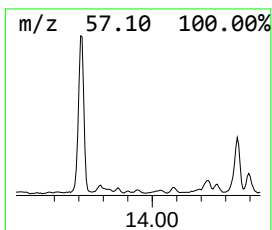
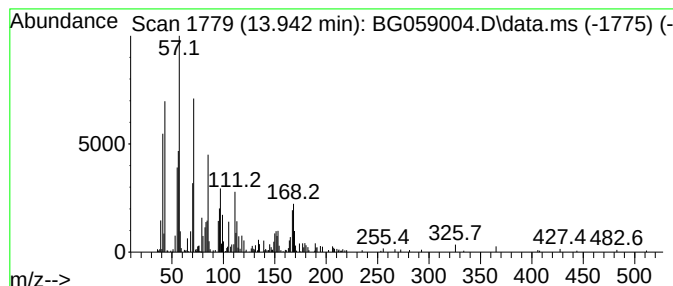
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.942	8.98 ng/ul	476246	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	55
2			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	50
3			1-Dodecanol, 2-hexyl-	270	C18H38O	110225-00-8	47
4			1-Dodecanol, 2-octyl-	298	C20H42O	005333-42-6	43
5			Heptane, 2,6-dimethyl-	128	C9H20	001072-05-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

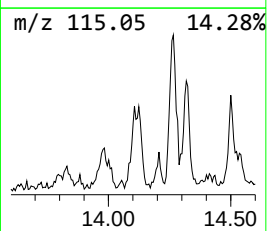
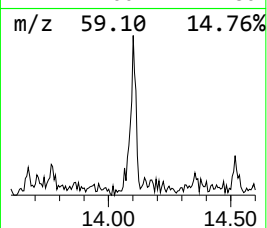
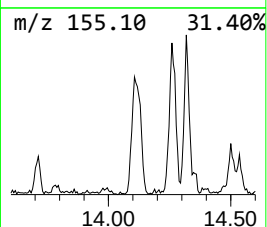
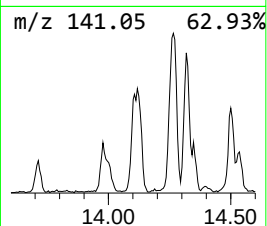
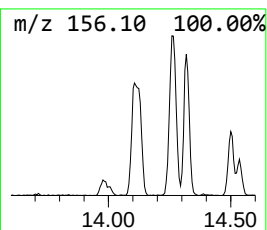
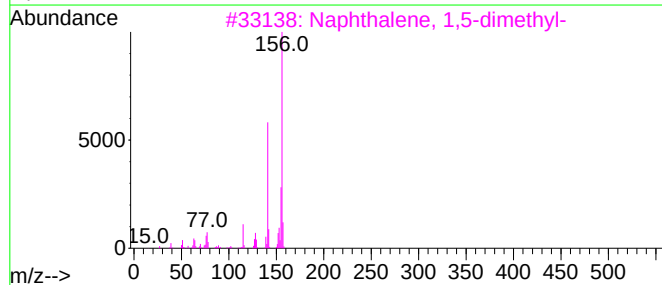
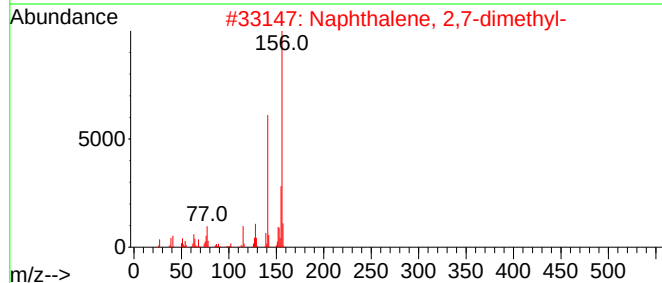
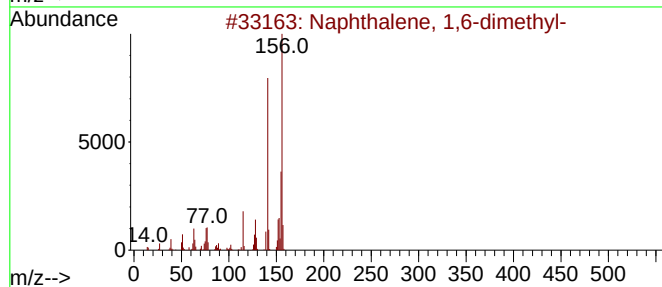
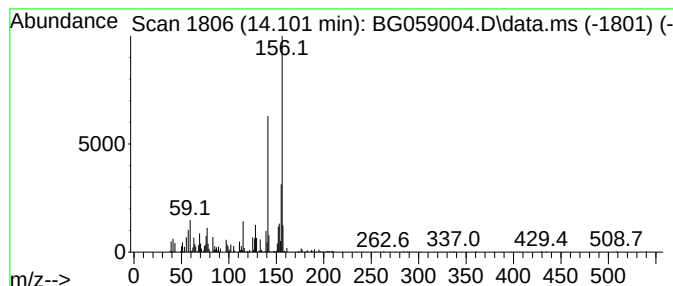
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 41 Naphthalene, 1,6-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.101	14.38 ng/ul	762976	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

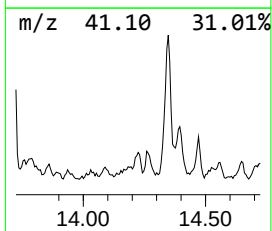
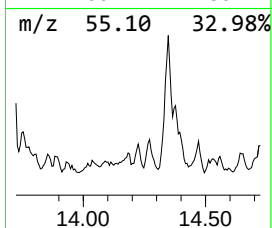
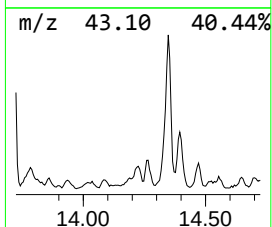
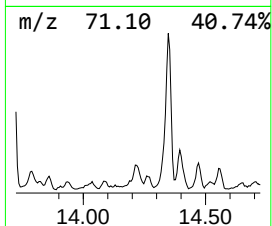
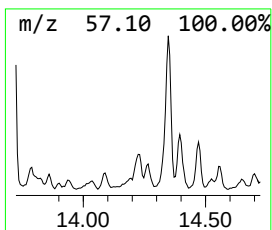
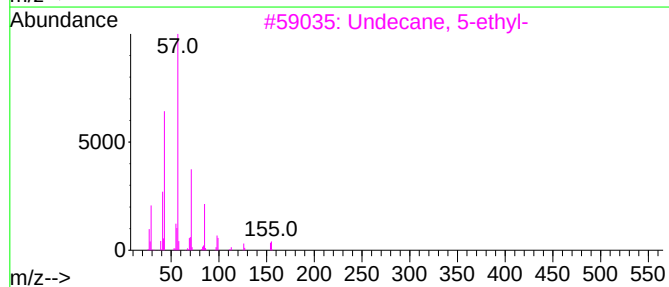
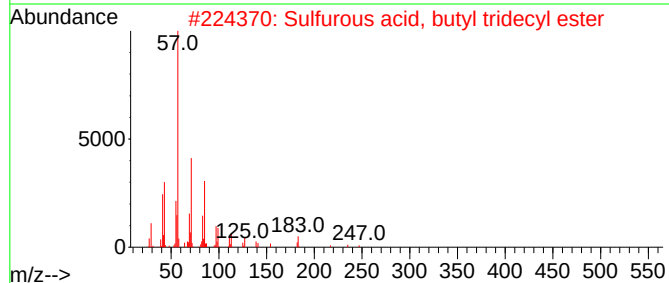
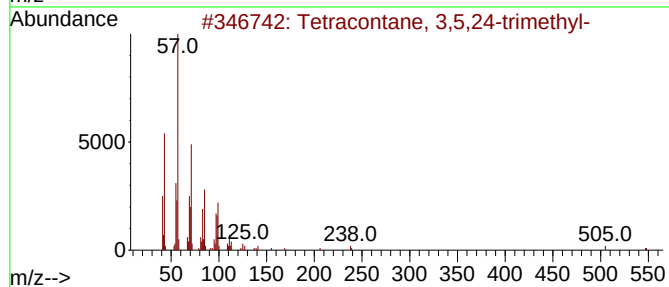
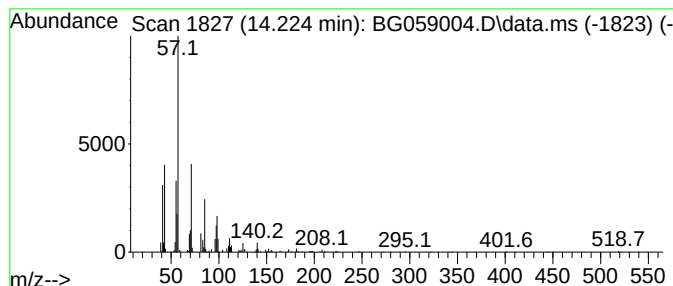
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 (DEL) Alkane: Straight-Chai... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.224	8.81 ng/ul	467479	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracontane, 3,5,24-trimethyl-	605	C43H88	055162-61-3	80
2			Sulfurous acid, butyl tridecyl e...	320	C17H36O3S	1000309-18-0	64
3			Undecane, 5-ethyl-	184	C13H28	017453-94-0	59
4			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	58
5			Nonyl tetradecyl ether	340	C23H48O	1000406-37-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

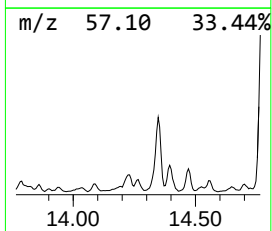
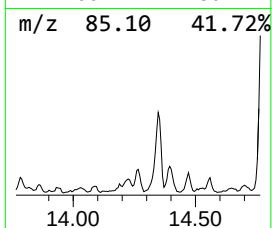
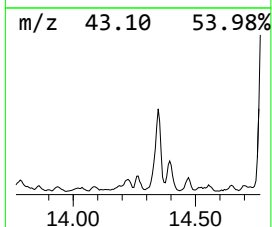
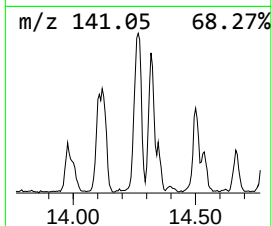
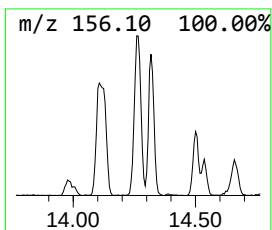
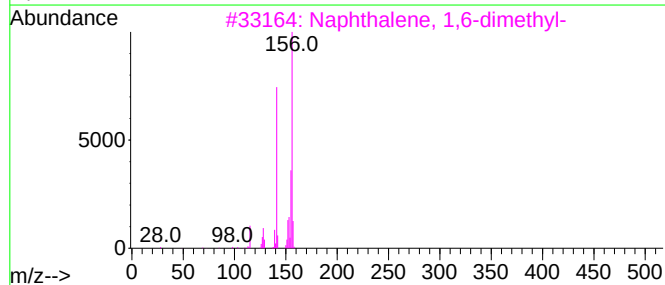
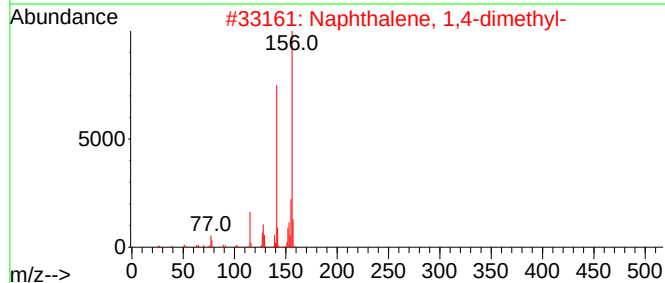
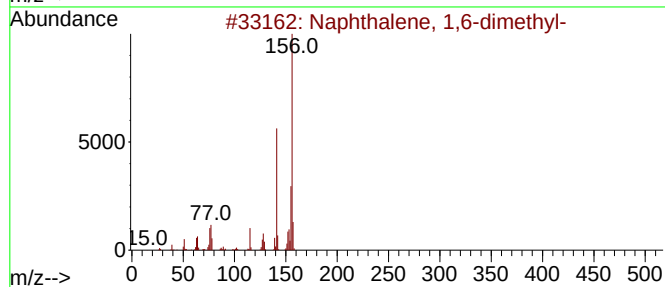
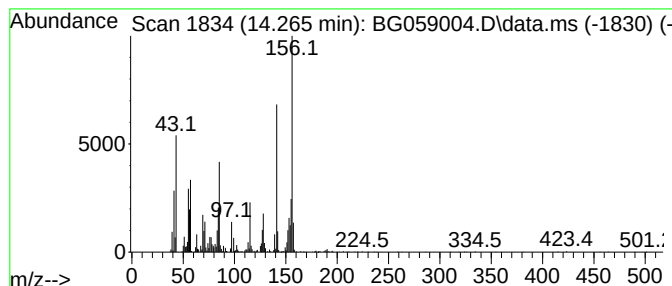
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 Naphthalene, 1,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.265	38.14 ng/ul	2023800	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	92
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	81
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	70
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

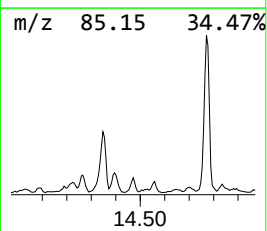
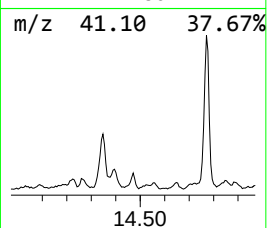
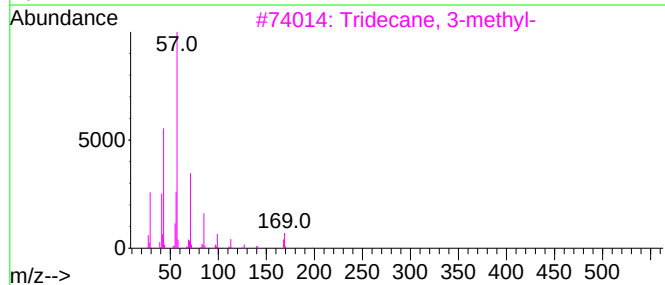
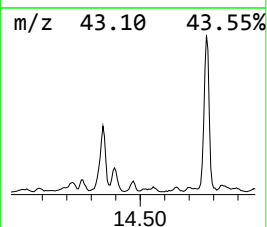
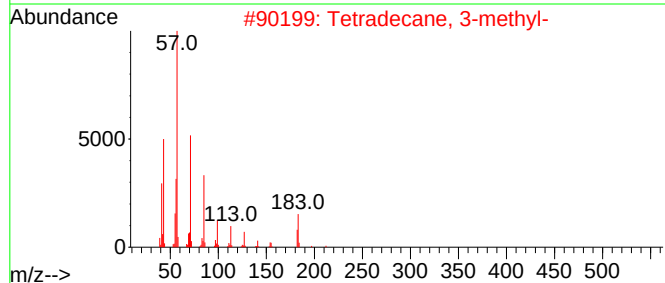
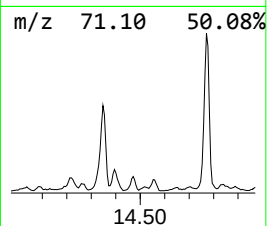
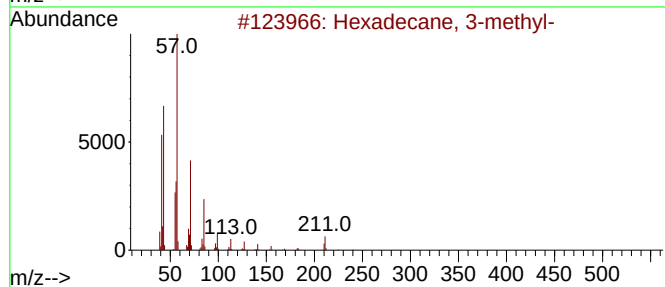
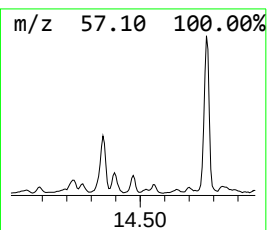
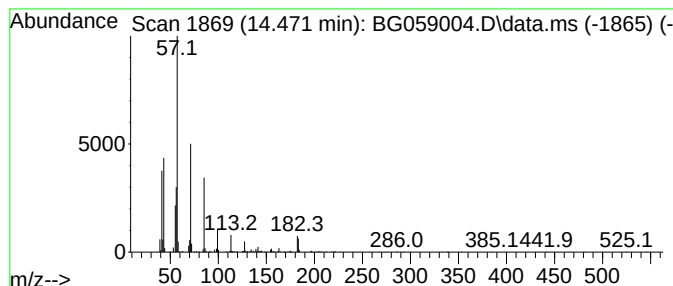
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.471	14.70 ng/ul	780141	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	90
2	Tetradecane, 3-methyl-	212	C15H32	018435-22-8	87
3	Tridecane, 3-methyl-	198	C14H30	006418-41-3	72
4	3,5-Dimethyldodecane	198	C14H30	107770-99-0	72
5	Tridecane, 7-hexyl-	268	C19H40	007225-66-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

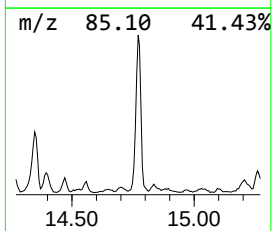
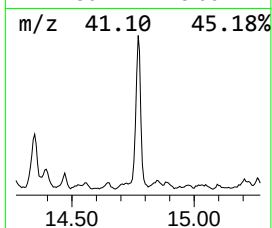
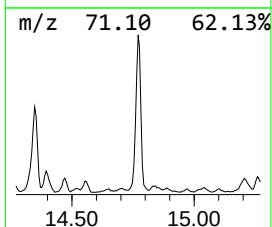
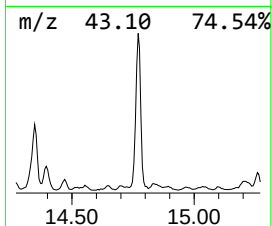
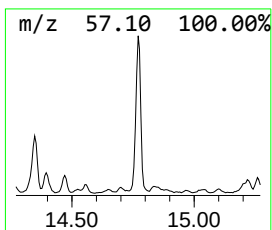
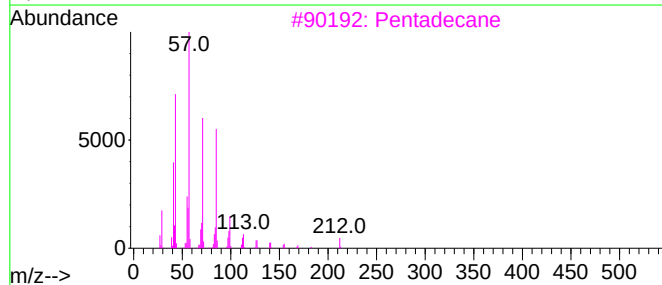
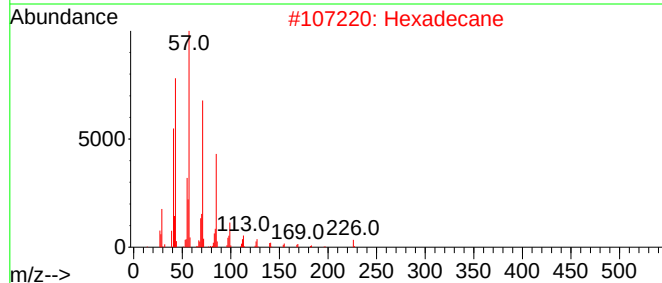
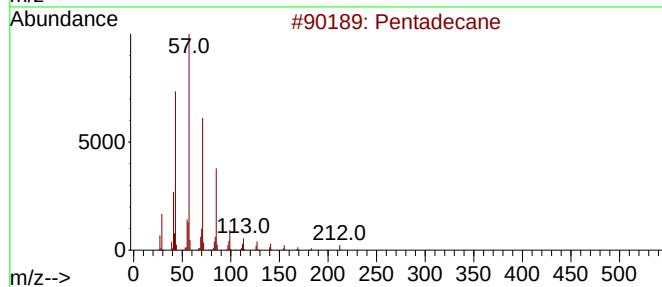
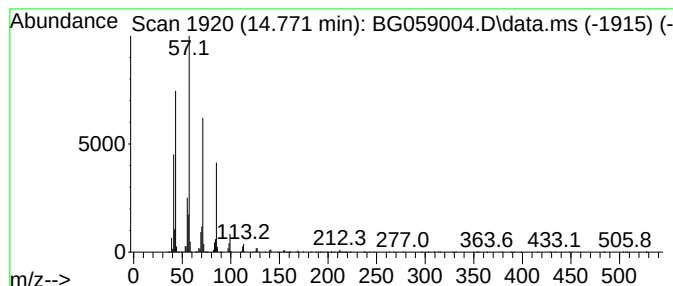
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.771	227.42 ng/ul	12067100	Acenaphthene-d10		14.947	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	91
2	Hexadecane		226	C16H34	000544-76-3	90
3	Pentadecane		212	C15H32	000629-62-9	87
4	Hexadecane		226	C16H34	000544-76-3	83
5	Tridecane		184	C13H28	000629-50-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

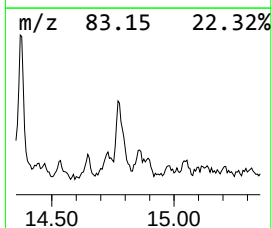
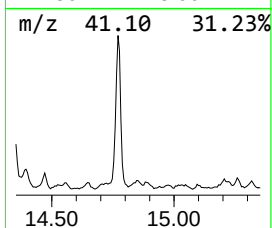
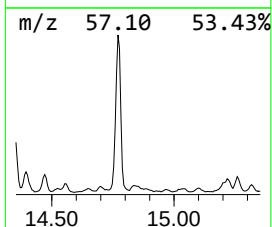
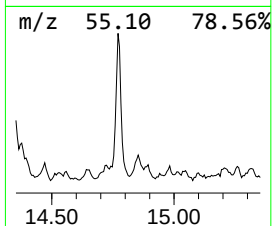
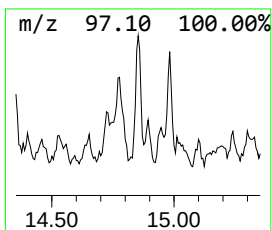
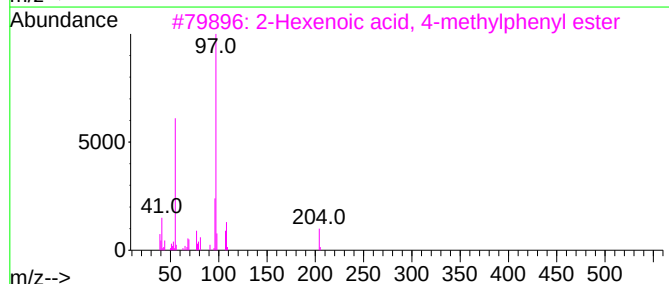
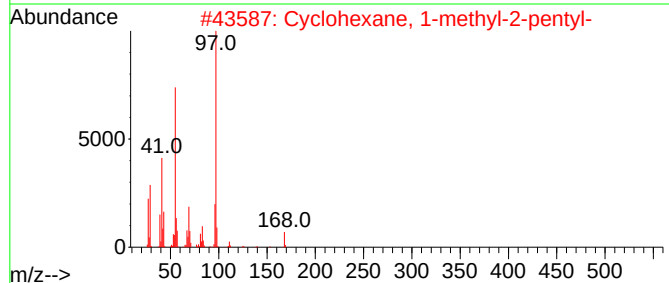
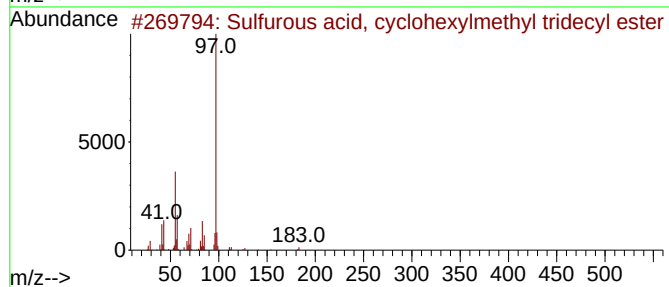
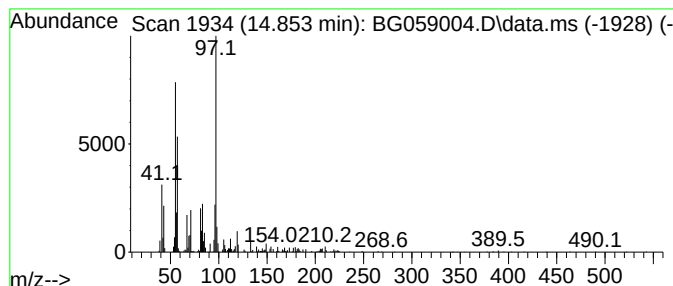
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 Sulfurous acid, cyclohexylm... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.853	18.30 ng/ul	970926	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, cyclohexylmethyl...	360	C20H40O3S	1000309-22-1	64
2	Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	59
3	2-Hexenoic acid, 4-methylphenyl ...	204	C13H16O2	069687-91-8	58
4	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	53
5	Sulfurous acid, cyclohexylmethyl...	234	C11H22O3S	1000309-21-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

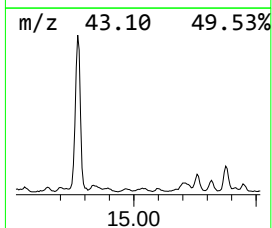
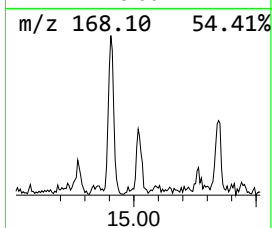
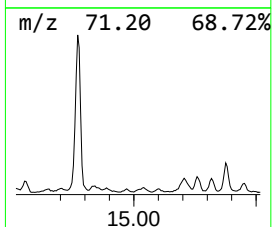
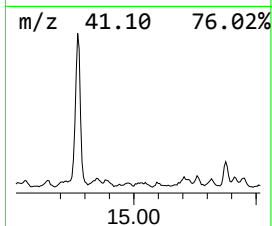
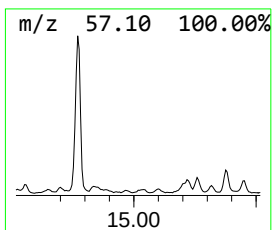
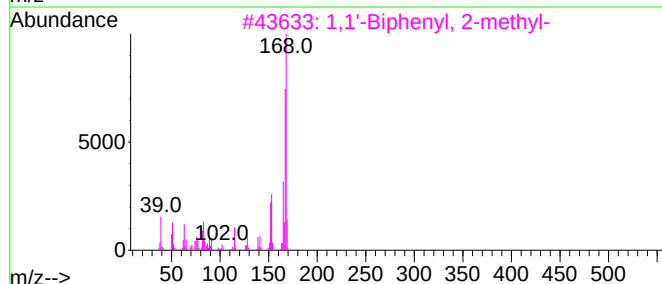
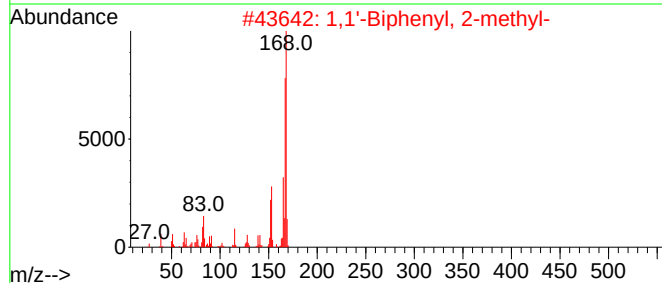
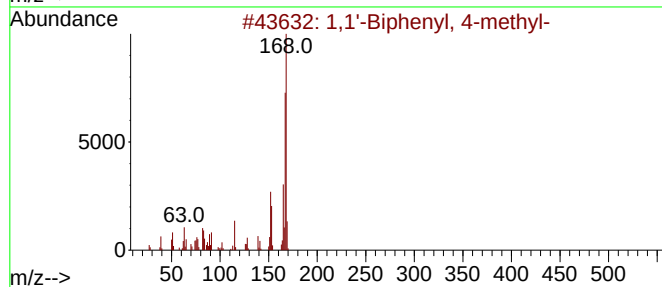
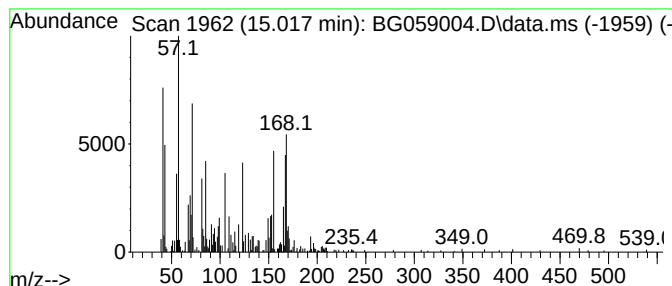
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 unknown-06 Concentration Rank 64

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.017	5.60 ng/ul	297383	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 4-methyl-			168	C13H12	000644-08-6	35
2	1,1'-Biphenyl, 2-methyl-			168	C13H12	000643-58-3	35
3	1,1'-Biphenyl, 2-methyl-			168	C13H12	000643-58-3	35
4	1,1'-Biphenyl, 3-methyl-			168	C13H12	000643-93-6	20
5	Tetradecane, 1-chloro-			232	C14H29Cl	002425-54-9	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

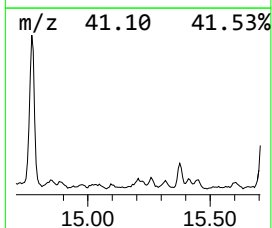
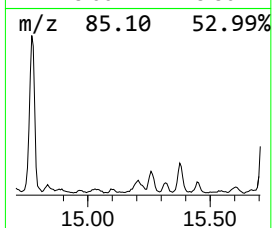
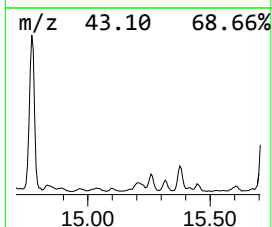
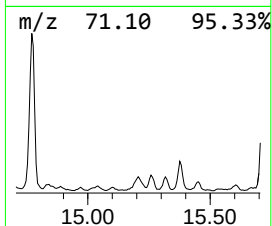
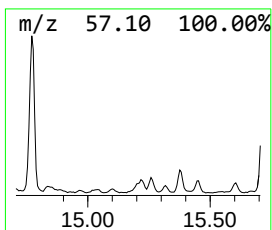
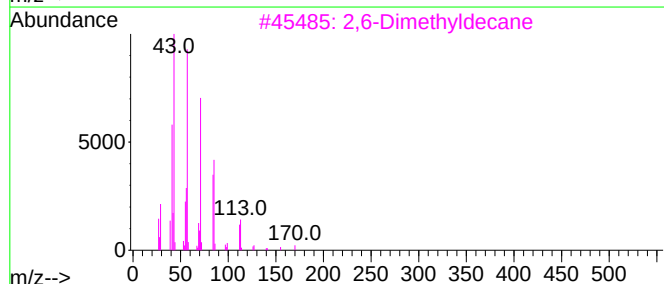
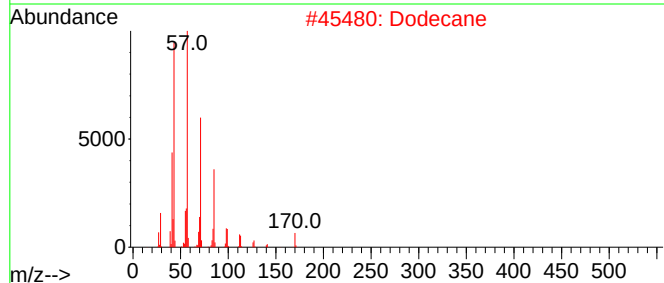
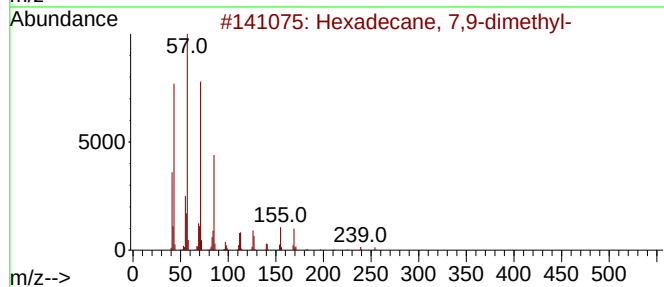
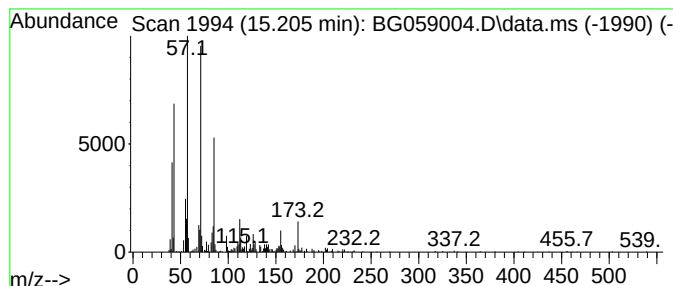
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.205	13.78 ng/ul	731422	Acenaphthene-d10		14.947	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane, 7,9-dimethyl-		254	C18H38	021164-95-4	80
2	Dodecane		170	C12H26	000112-40-3	80
3	2,6-Dimethyldecane		170	C12H26	013150-81-7	74
4	Tridecane, 5-propyl-		226	C16H34	055045-11-9	64
5	Tetradecane		198	C14H30	000629-59-4	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

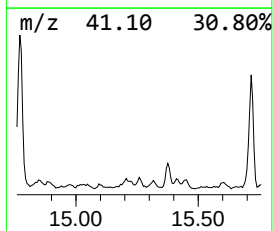
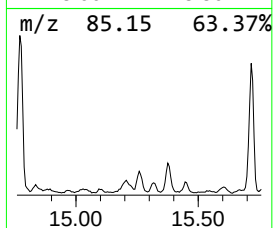
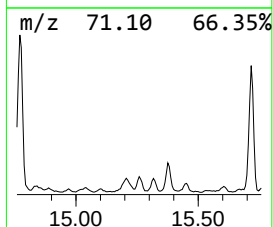
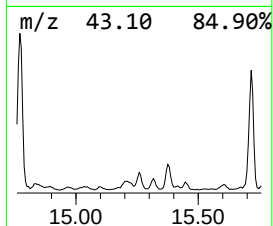
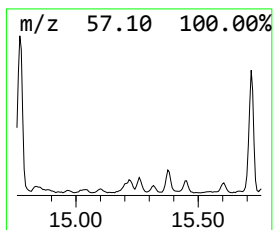
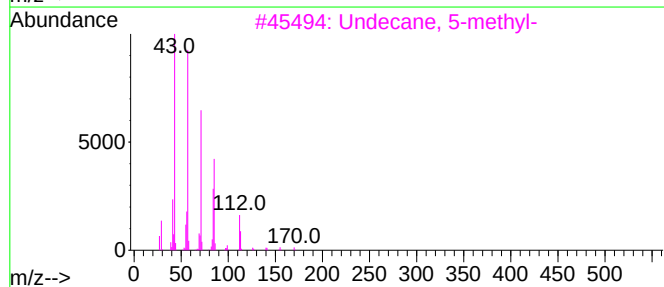
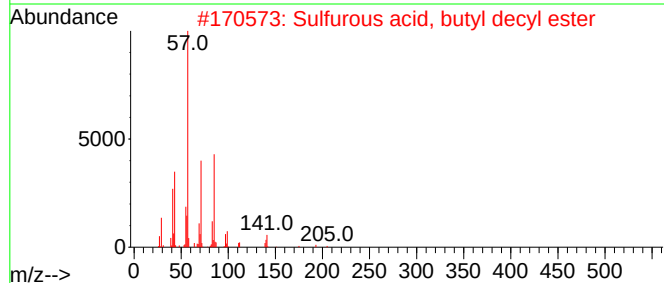
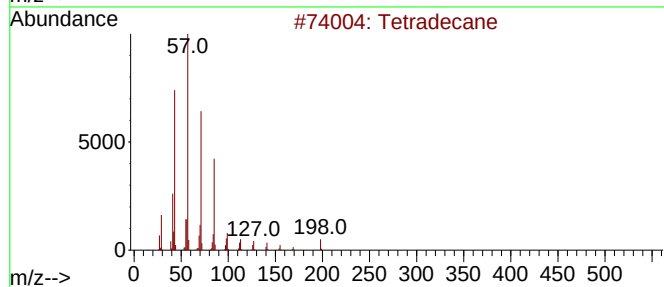
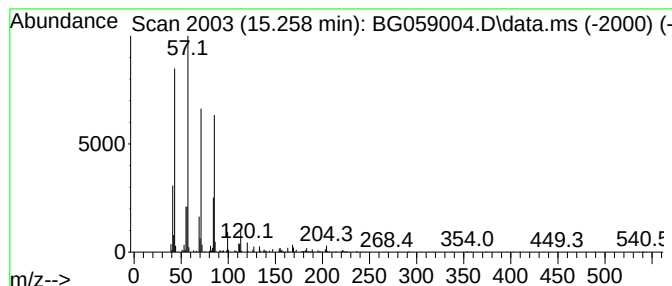
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.258	18.73 ng/ul	994079	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	72
2	Sulfurous acid, butyl decyl ester	278	C14H30O3S	1000309-17-7	59
3	Undecane, 5-methyl-	170	C12H26	001632-70-8	59
4	Decane, 2-methyl-	156	C11H24	006975-98-0	58
5	Tridecane, 5-methyl-	198	C14H30	025117-31-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

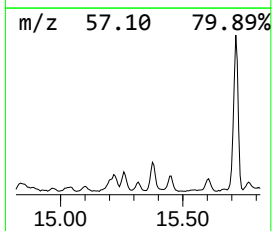
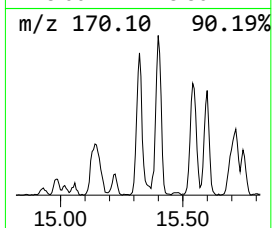
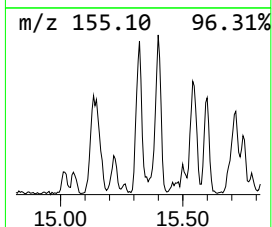
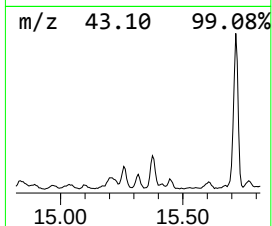
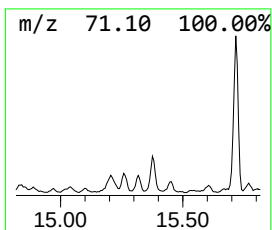
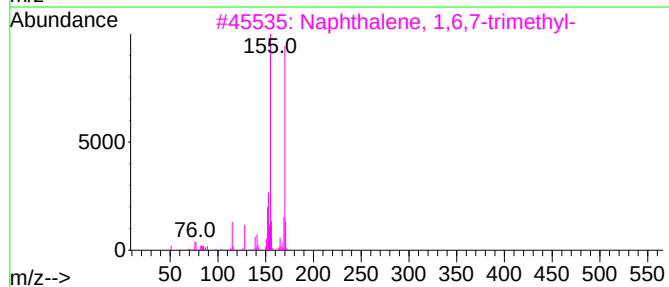
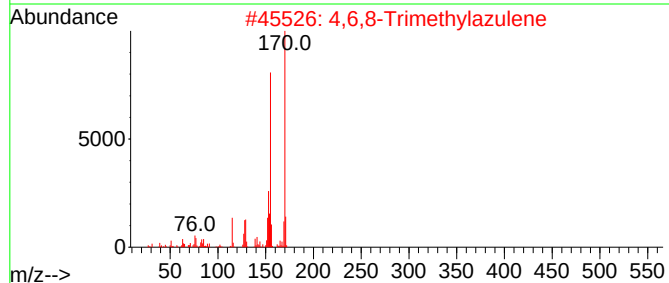
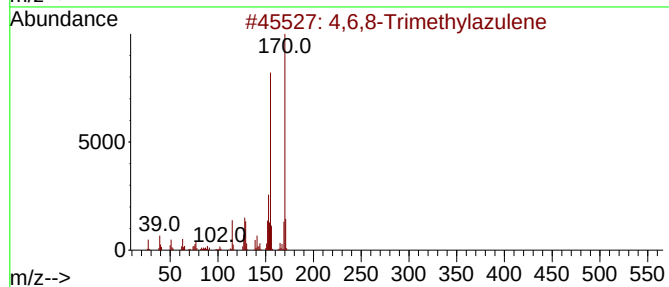
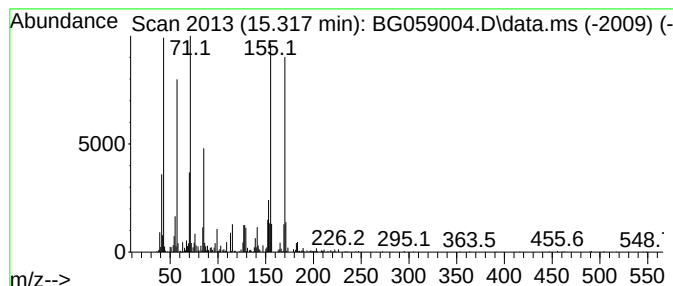
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 4,6,8-Trimethylazulene Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.317	22.43 ng/ul	1190380	Acenaphthene-d10	14.947

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	98
2		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	96
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
4		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90
5		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

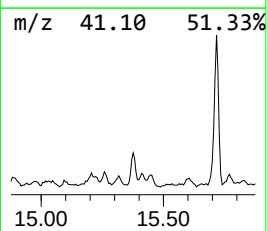
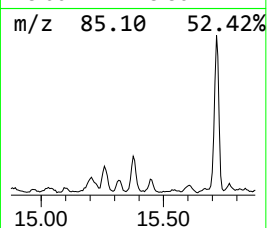
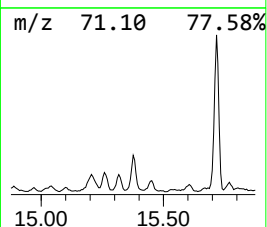
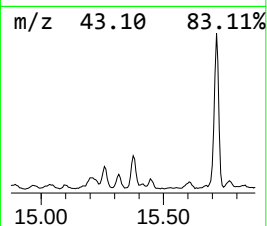
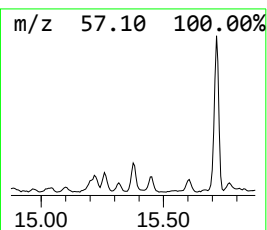
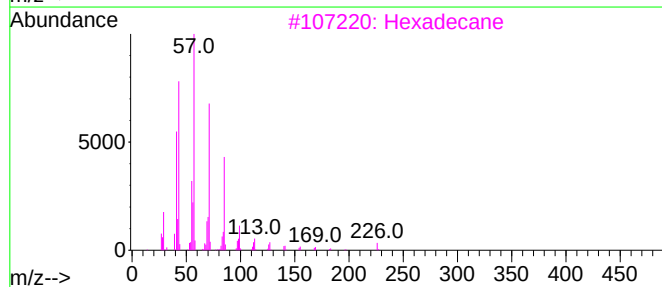
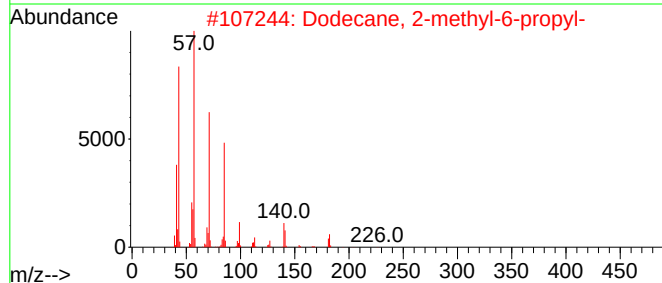
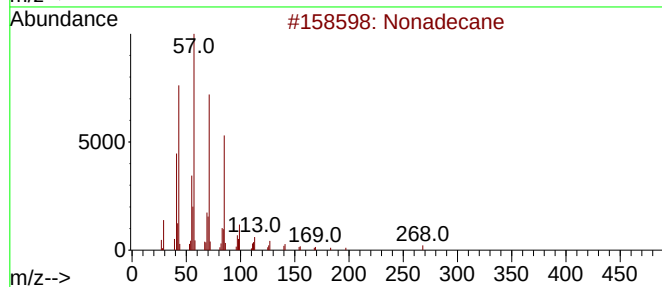
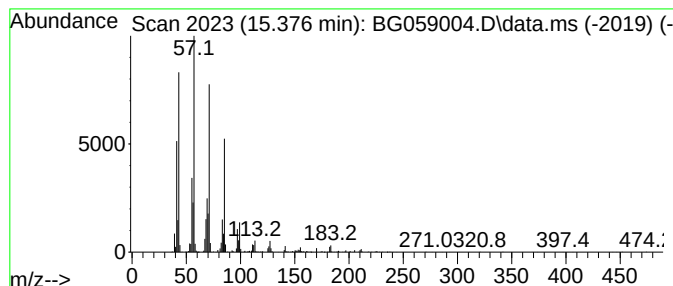
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.376	39.33 ng/ul	2086830	Acenaphthene-d10	14.947

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	91
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
3		Hexadecane	226	C16H34	000544-76-3	87
4		Dodecane	170	C12H26	000112-40-3	87
5		Dodecane	170	C12H26	000112-40-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

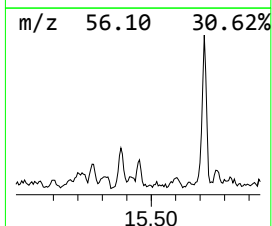
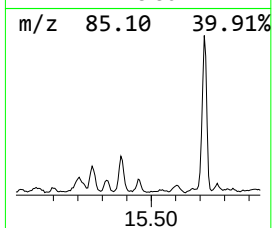
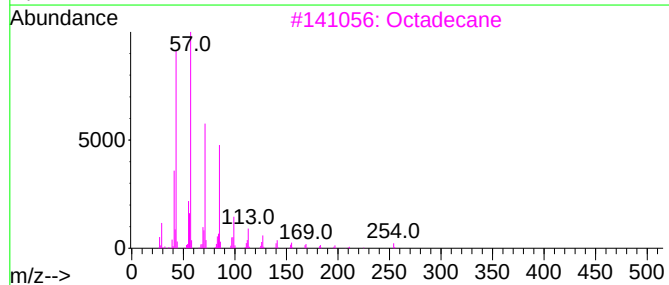
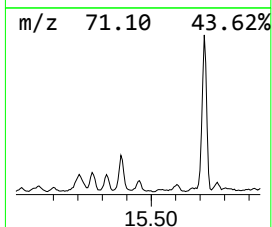
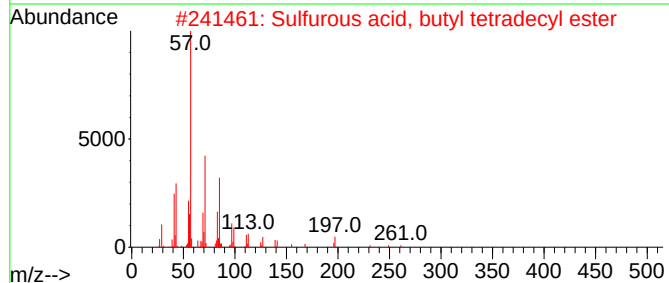
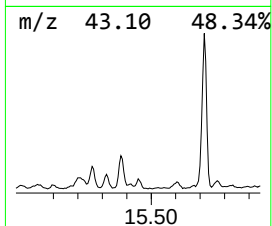
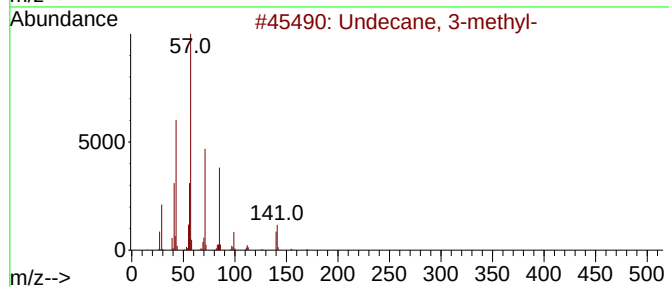
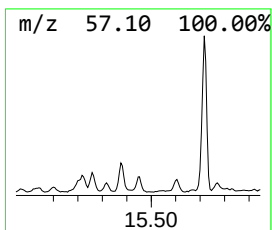
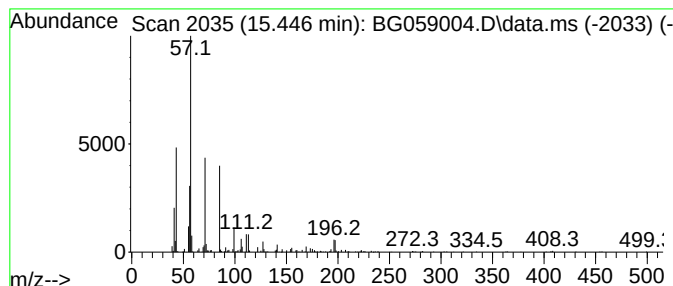
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.446	15.48 ng/ul	821256	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3-methyl-	170	C12H26	001002-43-3	72
2	Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	72
3	Octadecane	254	C18H38	000593-45-3	64
4	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	64
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

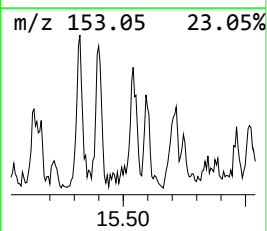
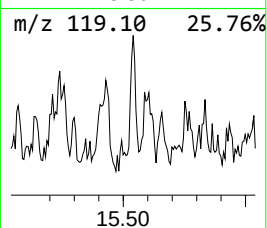
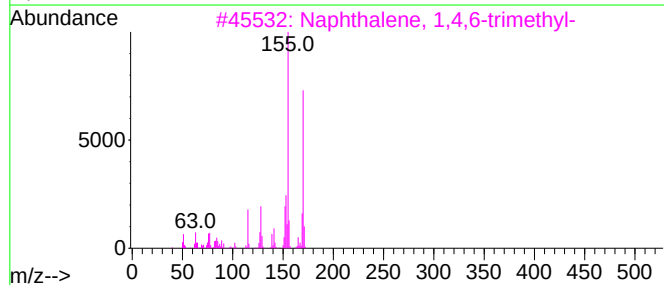
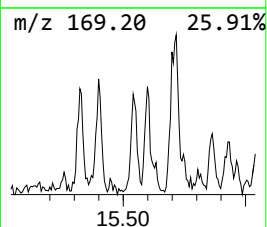
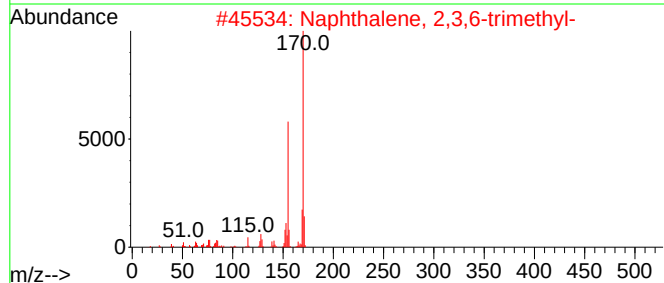
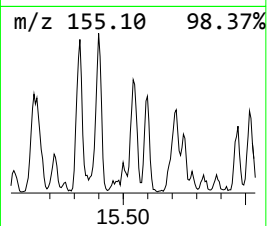
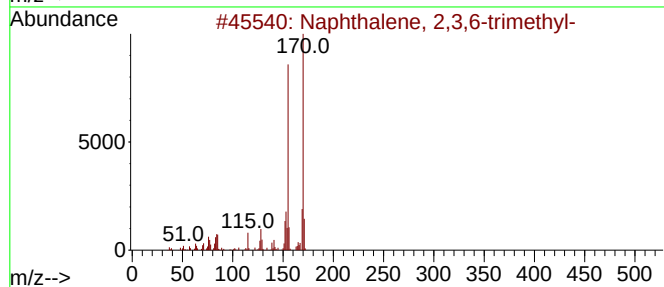
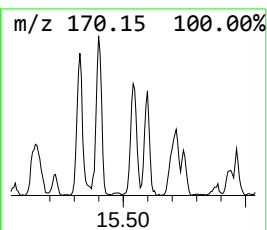
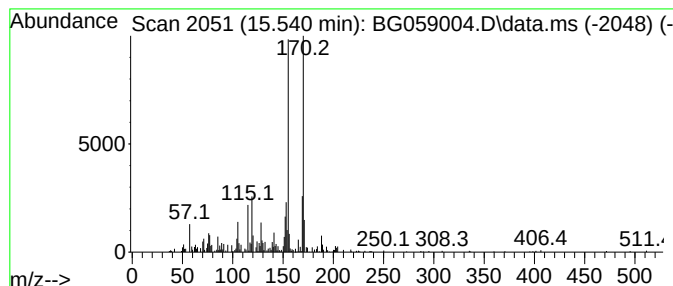
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 Naphthalene, 2,3,6-trimethyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.540	10.29 ng/ul	545971	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	91
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
5			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

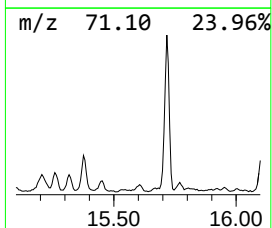
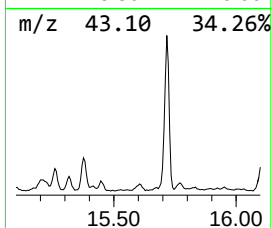
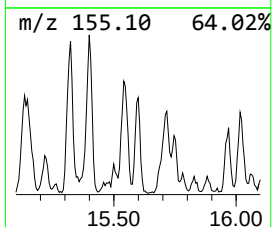
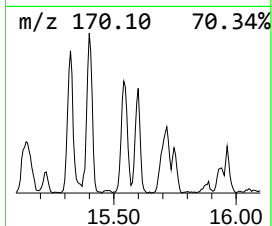
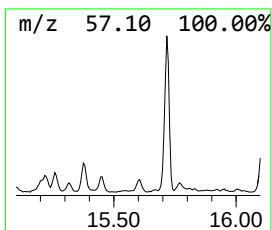
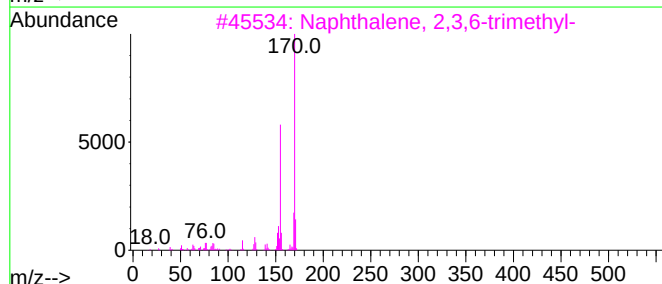
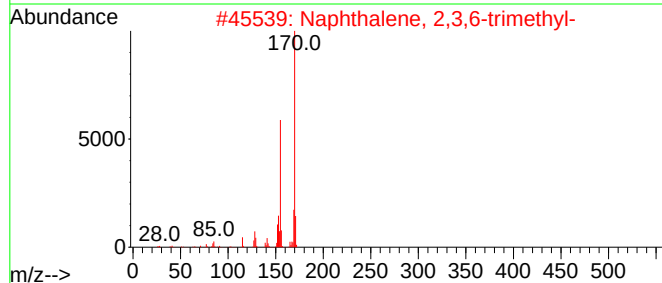
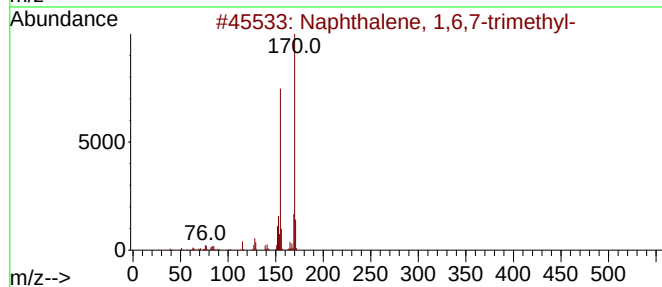
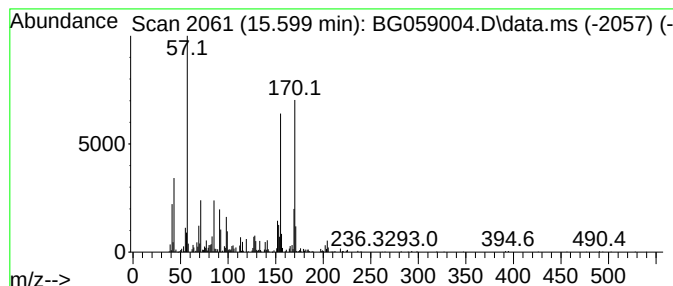
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 Naphthalene, 1,6,7-trimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.599	25.04 ng/ul	1328830	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
2	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	89
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

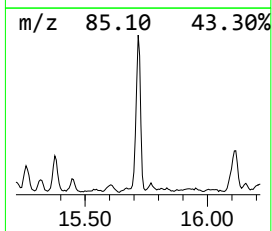
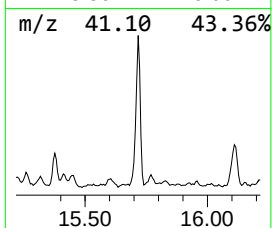
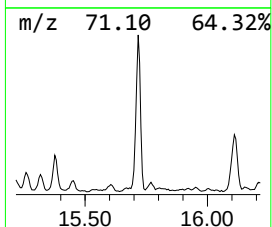
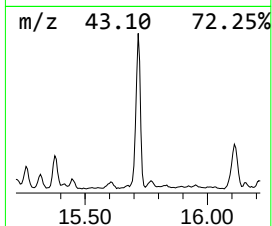
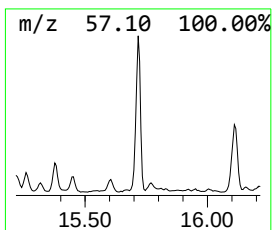
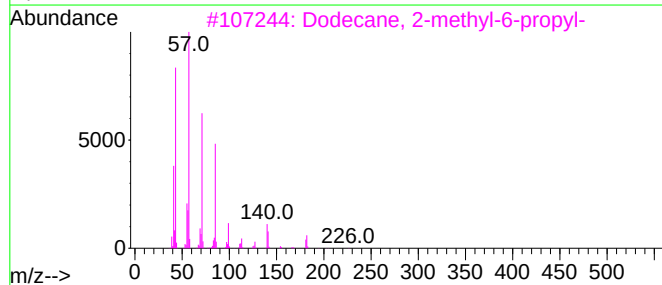
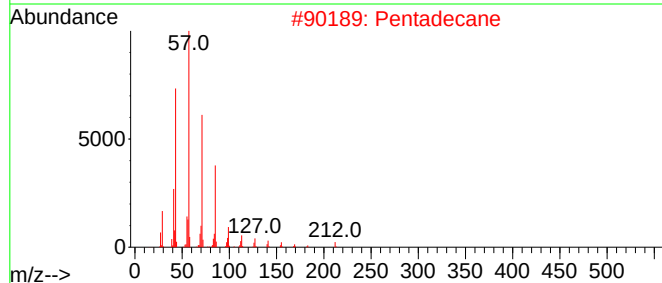
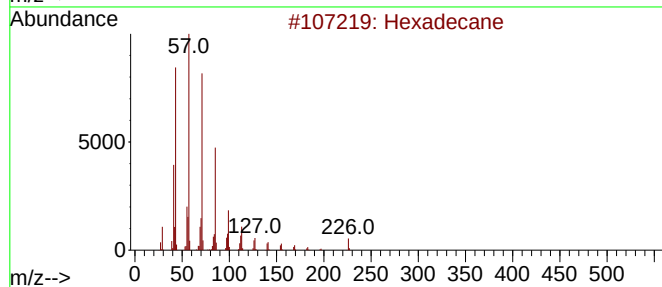
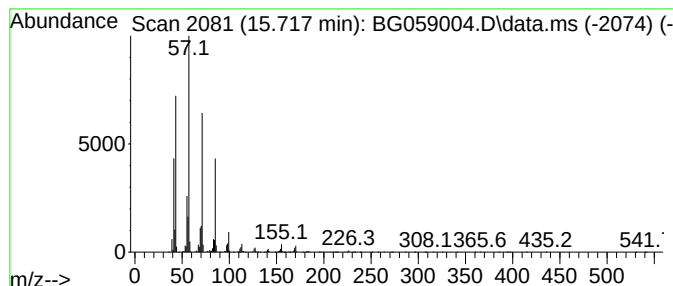
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.717	169.08 ng/ul	8971520	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Pentadecane	212	C15H32	000629-62-9	90
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
4	Dodecane, 4-methyl-	184	C13H28	006117-97-1	81
5	Dodecane, 2-methyl-	184	C13H28	001560-97-0	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

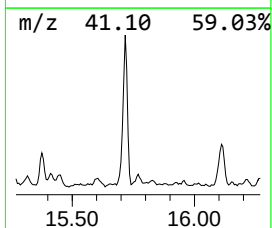
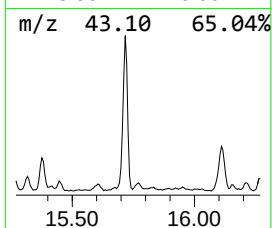
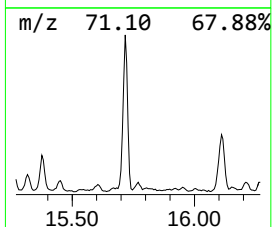
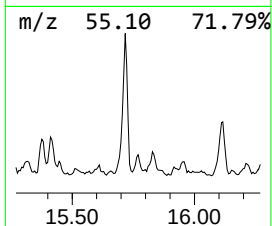
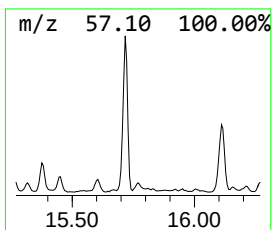
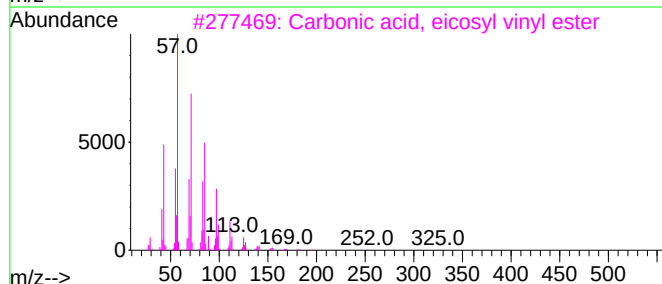
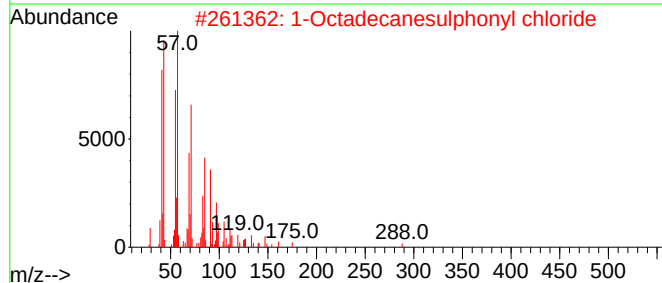
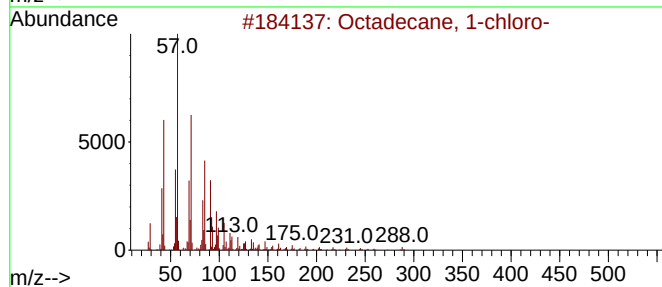
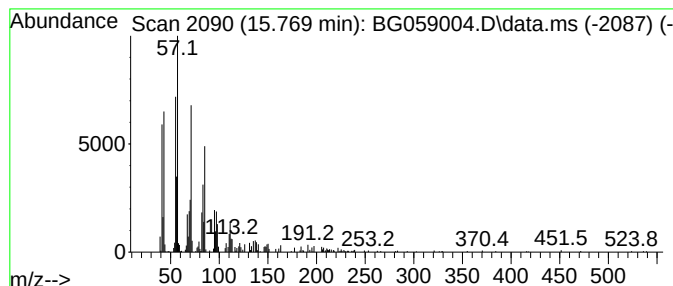
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 Octadecane, 1-chloro- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.770	11.81 ng/ul	626493	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	64
2			1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	64
3			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	64
4			Carbonic acid, decyl undecyl ester	356	C22H44O3	1000383-16-0	64
5			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

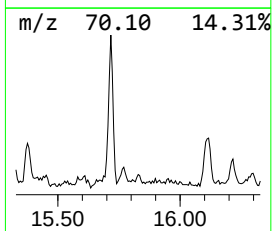
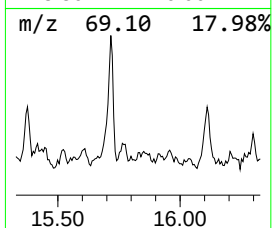
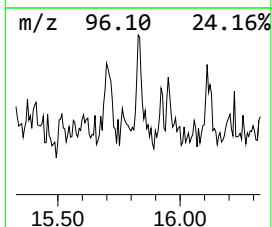
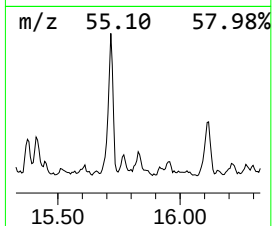
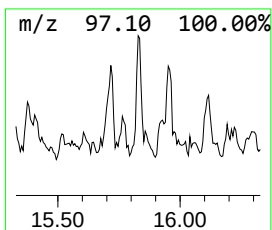
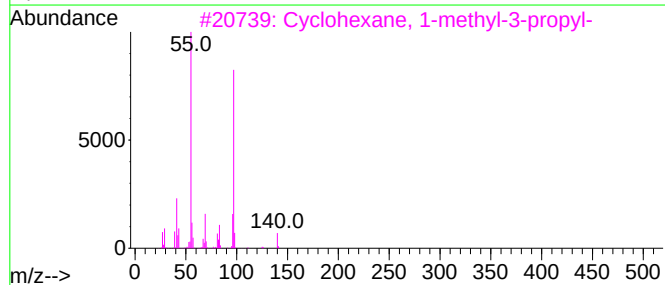
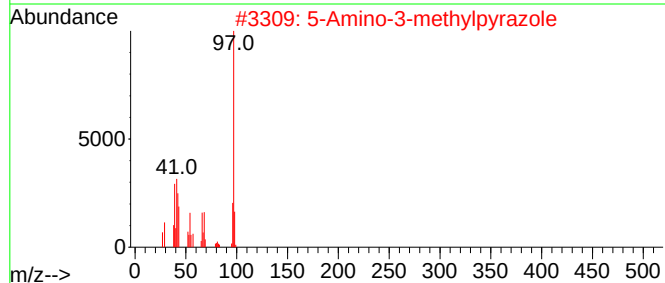
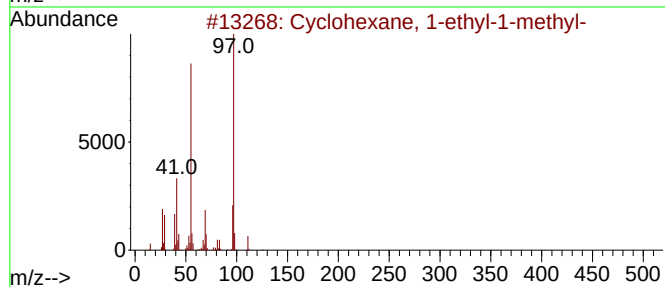
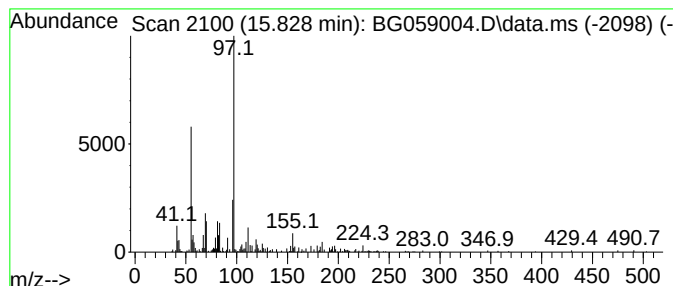
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 (DEL) Alkane: Cyclic15.828 Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.828	6.54 ng/ul	347271	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	72
2			5-Amino-3-methylpyrazole	97	C4H7N3	031230-17-8	59
3			Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	59
4			Sulfurous acid, cyclohexylmethyl...	262	C13H26O3S	1000309-21-5	50
5			Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

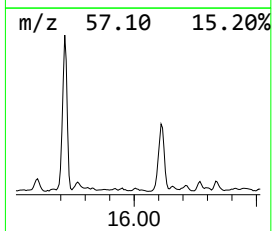
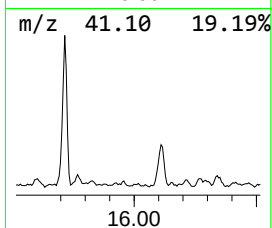
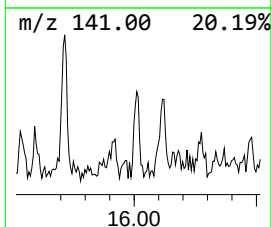
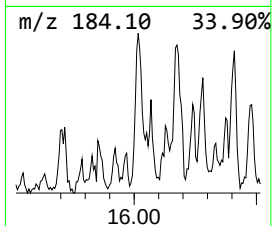
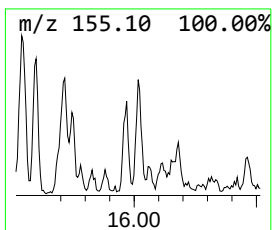
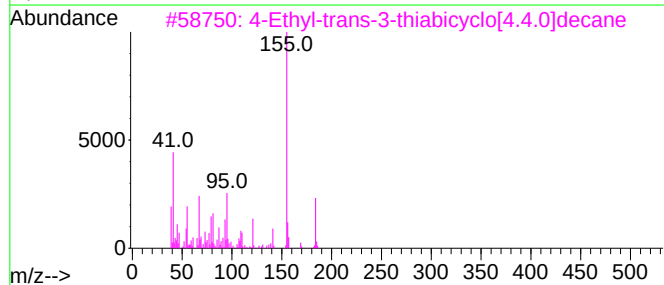
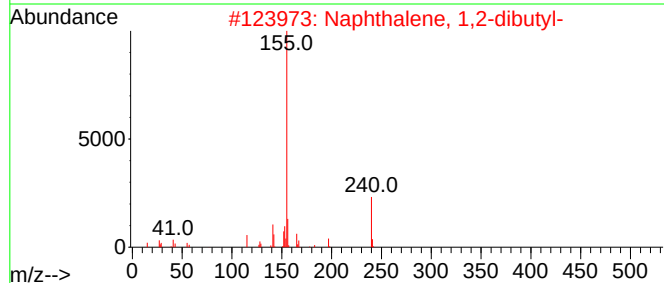
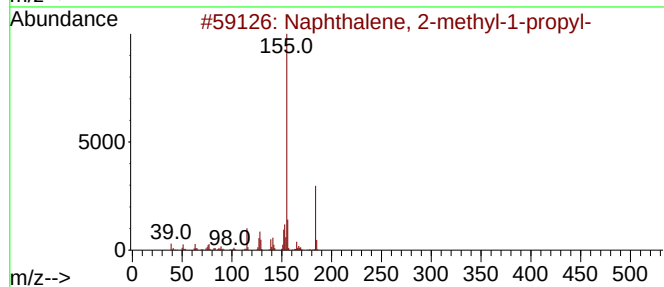
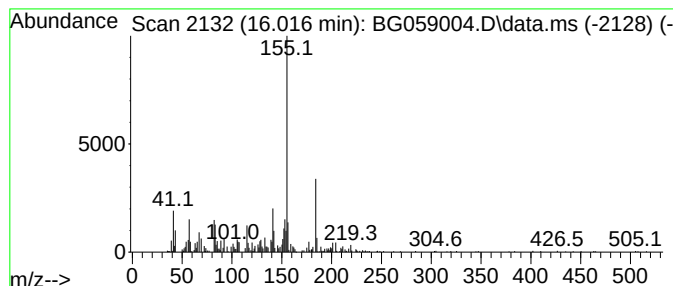
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 58 Naphthalene, 2-methyl-1-pro... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.016	10.78 ng/ul	571881	Acenaphthene-d10	14.947

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-1-propyl-	184	C14H16	054774-89-9	87
2			Naphthalene, 1,2-dibutyl-	240	C18H24	103385-90-6	53
3			4-Ethyl-trans-3-thiabicyclo[4.4....	184	C11H20S	1000215-94-2	53
4			1,2-Cyclohexanedicarboxylic acid...	354	C21H38O4	1000339-50-6	47
5			Benzamide, 2,6-difluoro-3-methyl...	227	C12H15F2NO	1000407-73-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

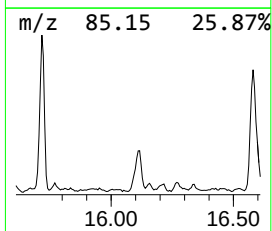
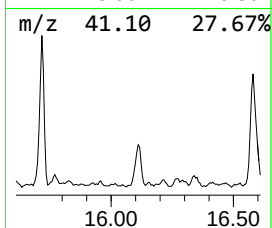
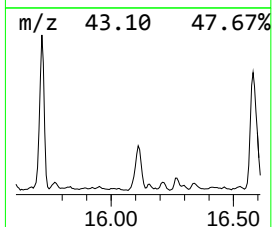
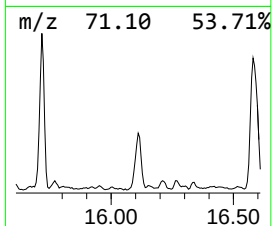
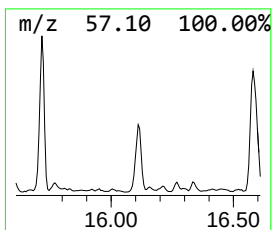
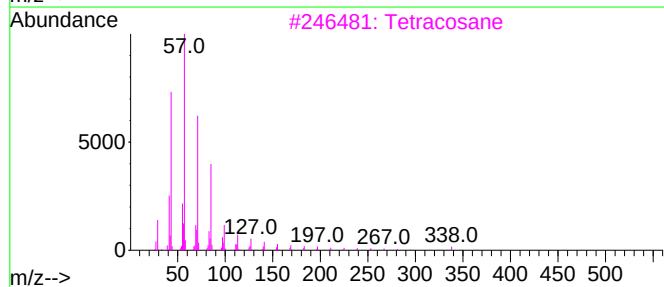
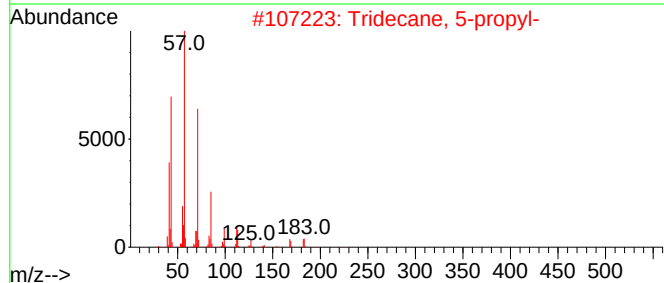
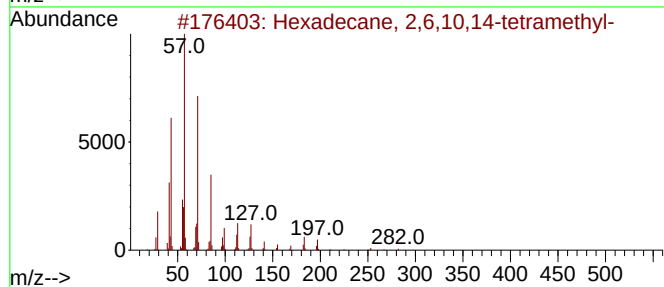
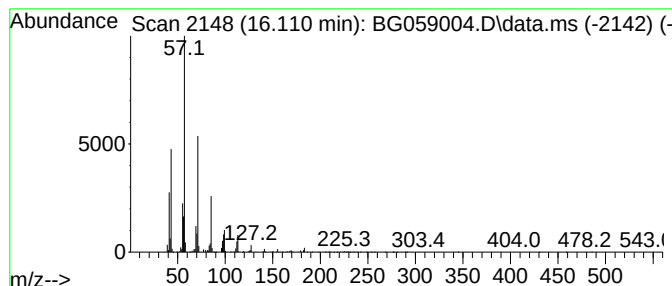
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.110	74.76 ng/ul	3966650	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	83
2	Tridecane, 5-propyl-	226	C16H34	055045-11-9	81
3	Tetracosane	338	C24H50	000646-31-1	80
4	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	72
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

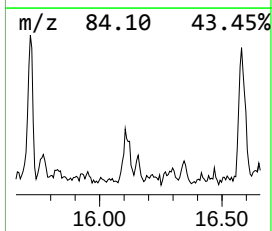
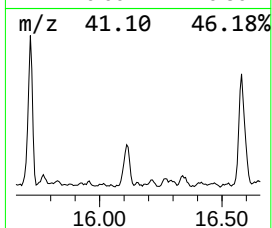
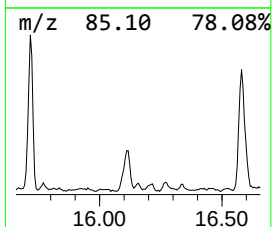
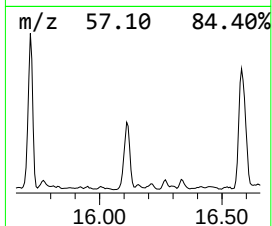
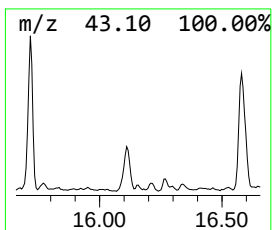
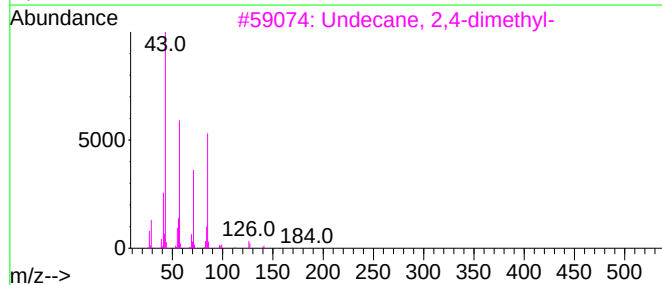
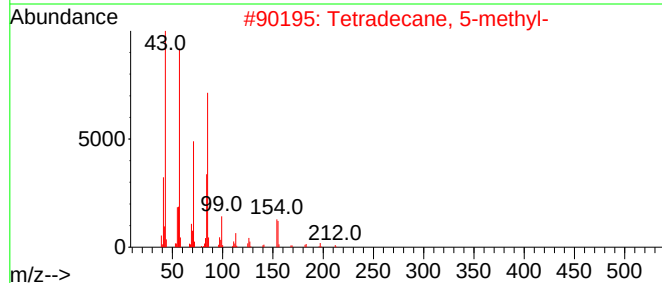
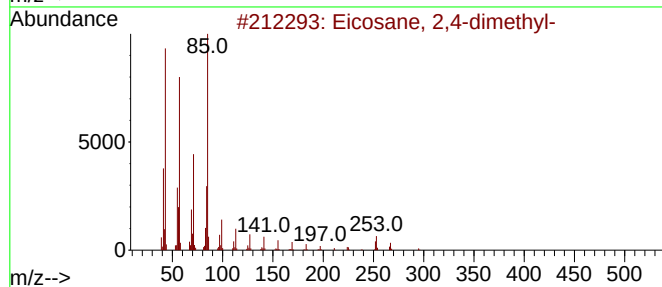
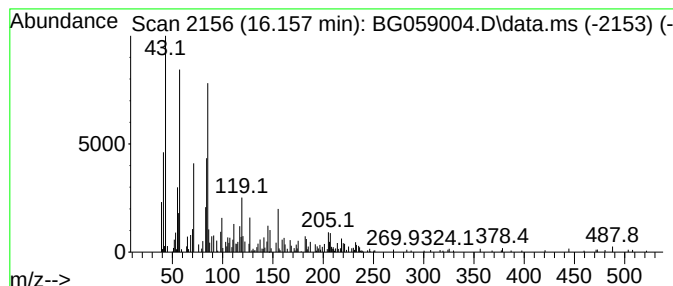
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.157	4.79 ng/ul	254208	Acenaphthene-d10		14.947	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane, 2,4-dimethyl-		310	C22H46	075163-98-3	50
2	Tetradecane, 5-methyl-		212	C15H32	025117-32-2	47
3	Undecane, 2,4-dimethyl-		184	C13H28	017312-80-0	47
4	tert-Butyldimethylsilyl nitrile		141	C7H15NSi	056522-24-8	46
5	Nonane, 5-methyl-		142	C10H22	015869-85-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

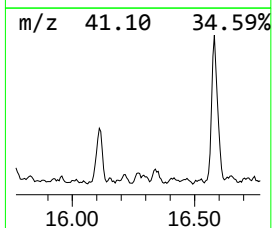
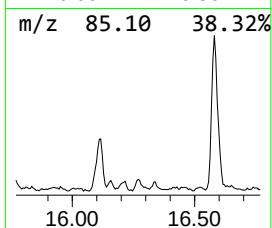
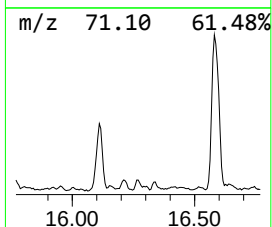
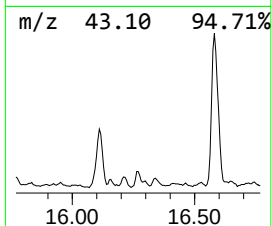
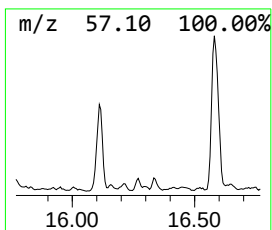
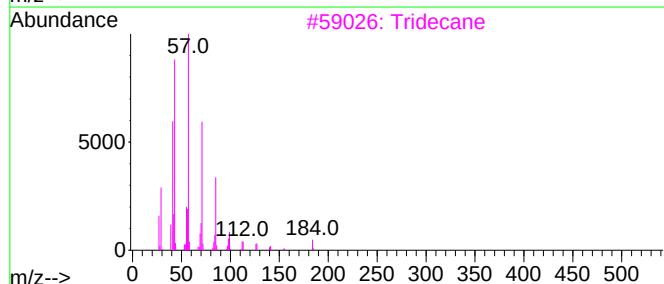
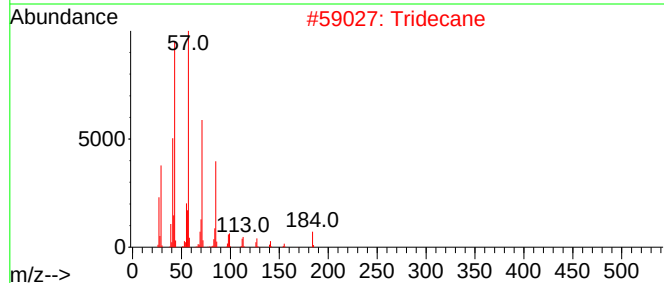
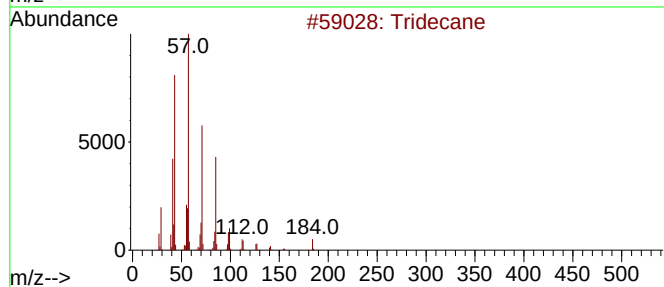
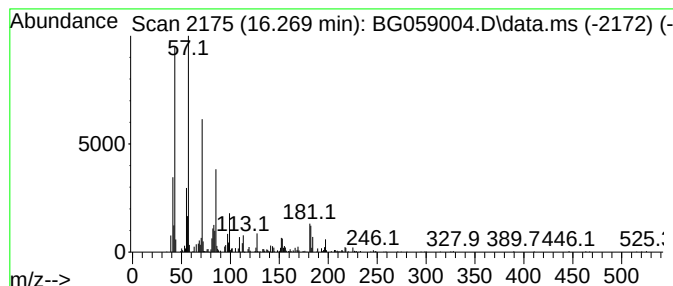
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.269	11.76 ng/ul	623905	Acenaphthene-d10	14.947	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	90
2	Tridecane	184	C13H28	000629-50-5	81
3	Tridecane	184	C13H28	000629-50-5	81
4	Tetracosane	338	C24H50	000646-31-1	72
5	Heneicosane	296	C21H44	000629-94-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

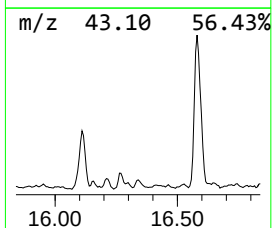
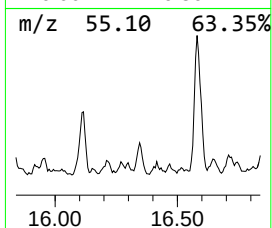
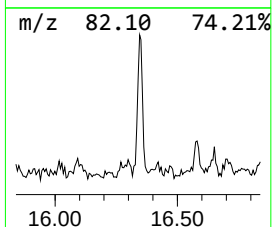
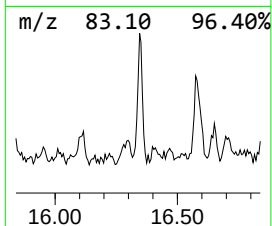
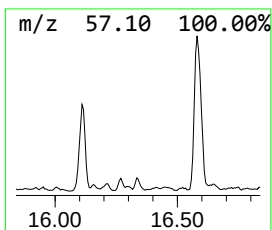
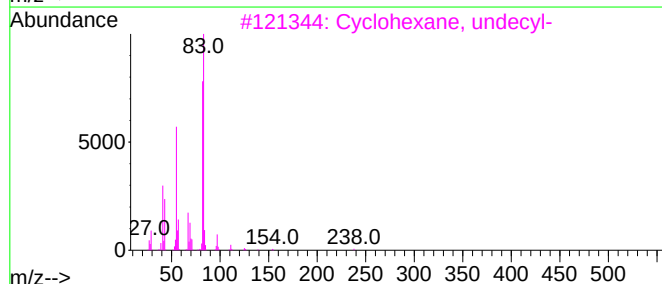
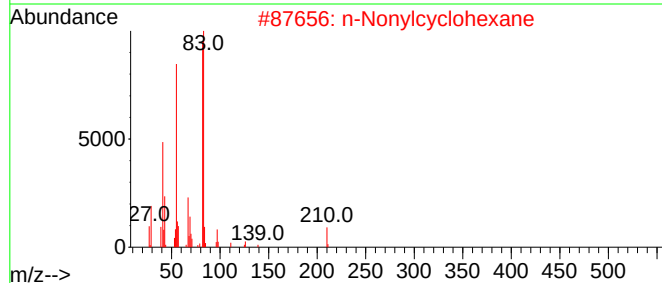
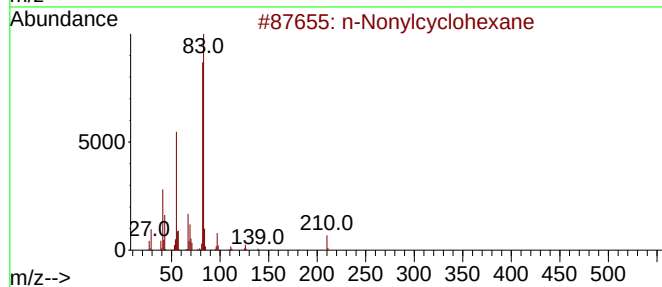
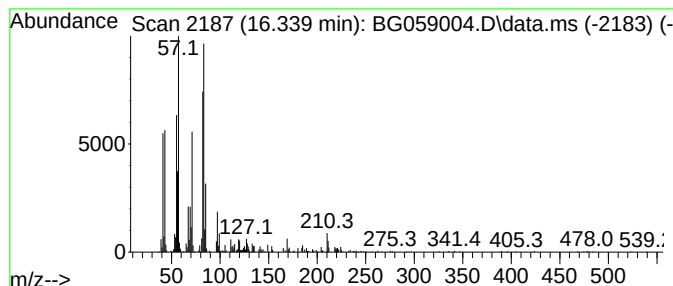
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 (DEL) Alkane: Cyclic16.339 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.339	17.11 ng/ul	969264	Phenanthrene-d10		17.697	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Nonylcyclohexane		210	C15H30	002883-02-5	58
2	n-Nonylcyclohexane		210	C15H30	002883-02-5	58
3	Cyclohexane, undecyl-		238	C17H34	054105-66-7	53
4	n-Tridecylcyclohexane		266	C19H38	006006-33-3	53
5	Cyclohexane, decyl-		224	C16H32	001795-16-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

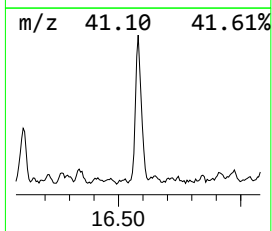
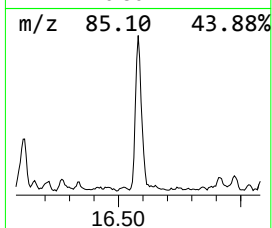
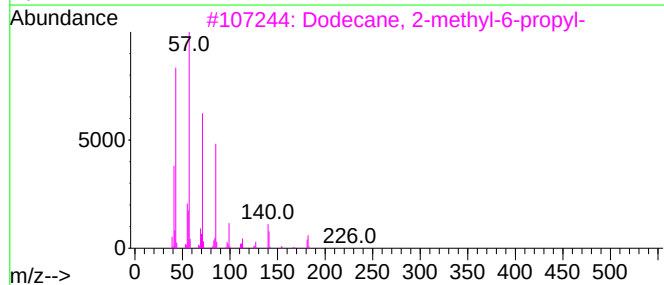
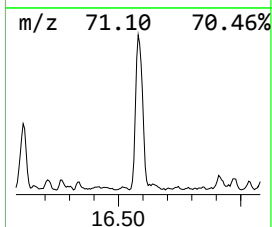
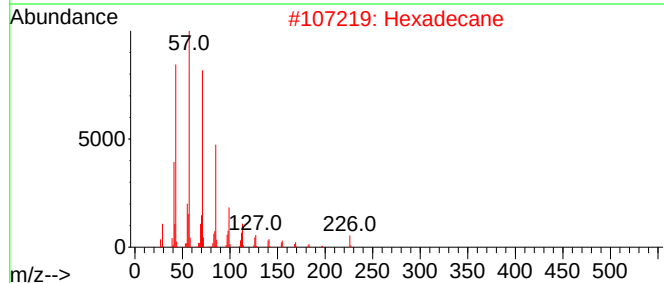
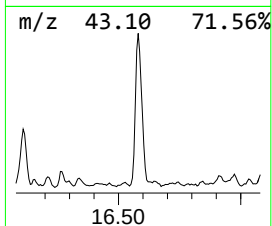
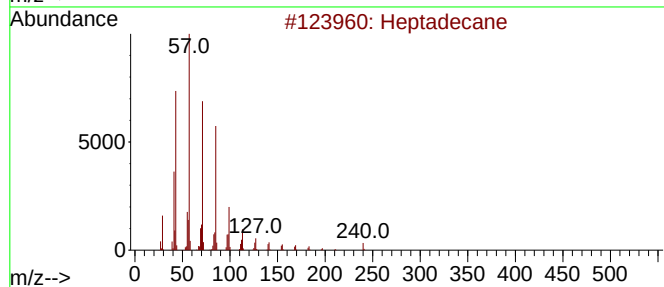
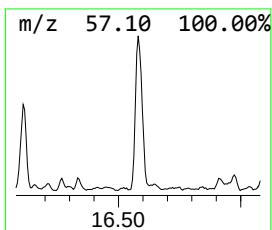
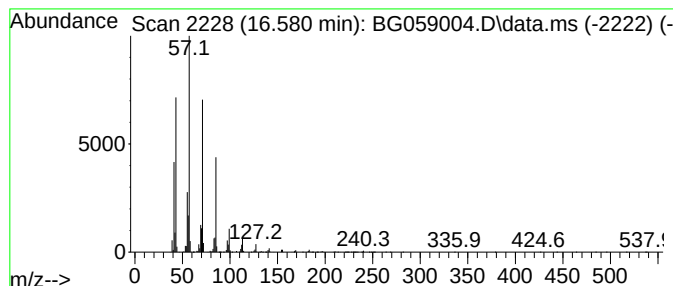
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.580	173.16 ng/ul	9807630	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	95
2			Hexadecane	226	C16H34	000544-76-3	94
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4			Tritetracontane	605	C43H88	007098-21-7	90
5			Pentadecane	212	C15H32	000629-62-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

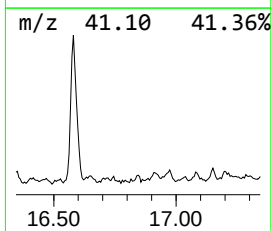
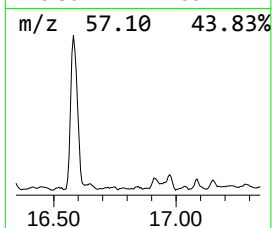
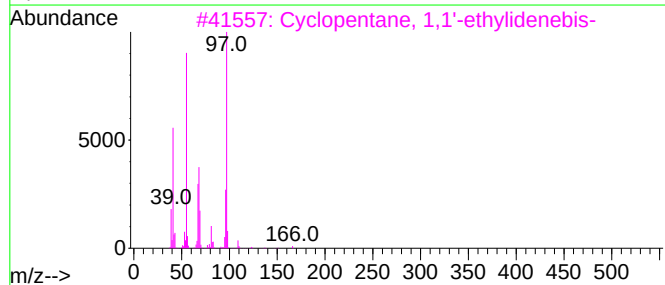
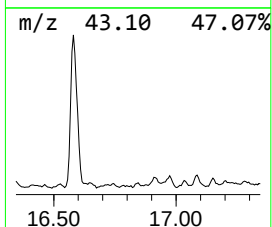
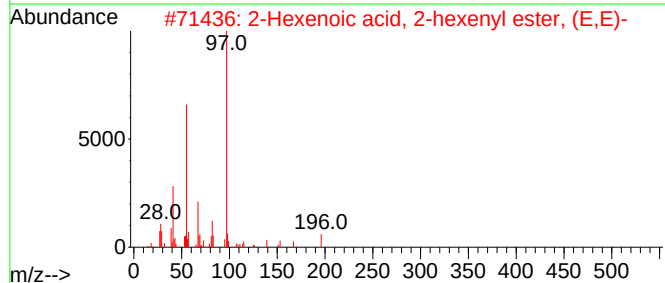
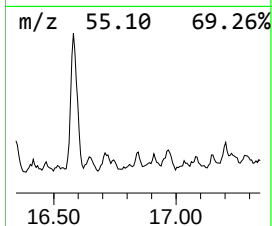
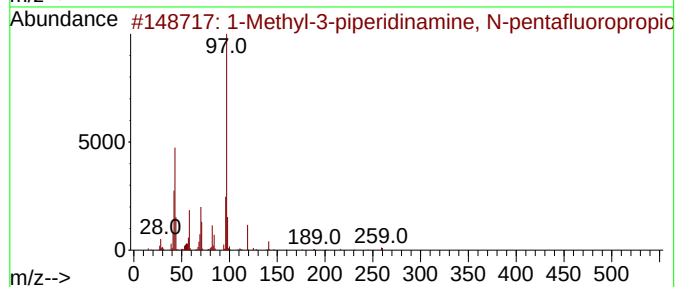
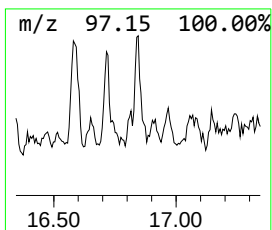
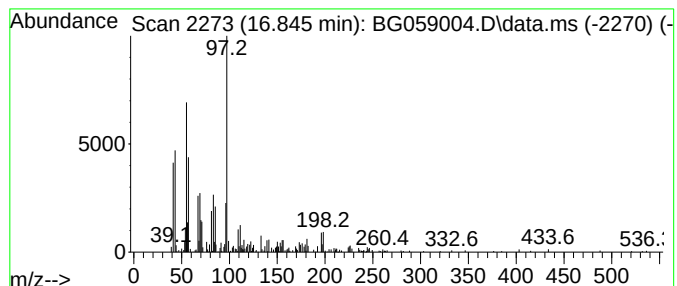
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 64 1-Methyl-3-piperidinamine, ... Concentration Rank 67

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.845	4.93 ng/ul	279438	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-3-piperidinamine, N-pen...	260	C9H13F5N2O	1963071-64-8	58
2			2-Hexenoic acid, 2-hexenyl ester...	196	C12H20O2	054845-28-2	53
3			Cyclopentane, 1,1'-ethylidenebis-	166	C12H22	004413-21-2	52
4			Succinic acid, cyclohexylmethyl ...	310	C18H30O4	1000391-41-3	50
5			Pyrollidine, 2,5-bis(imino)-	97	C4H7N3	000503-84-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

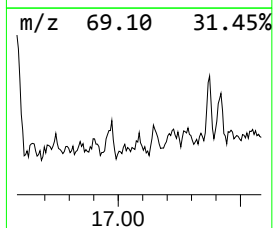
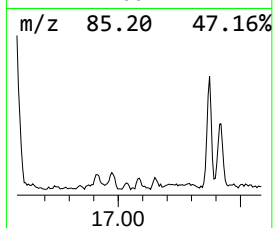
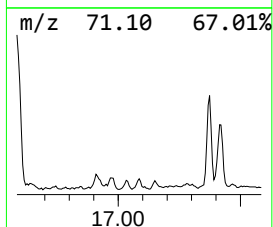
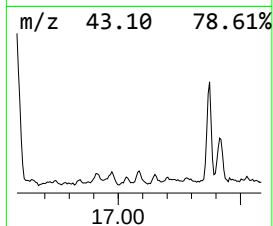
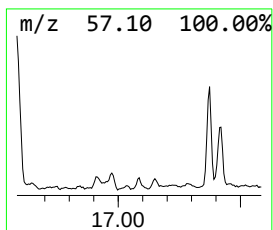
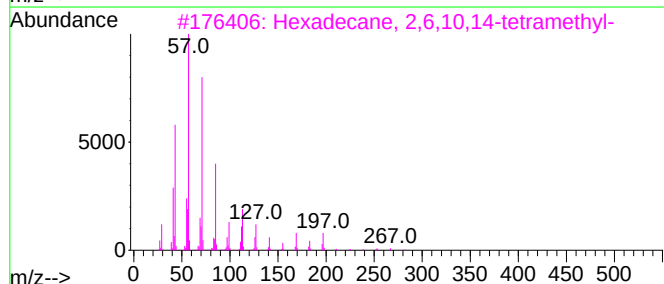
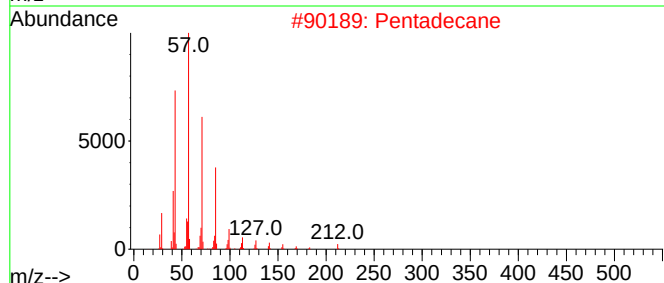
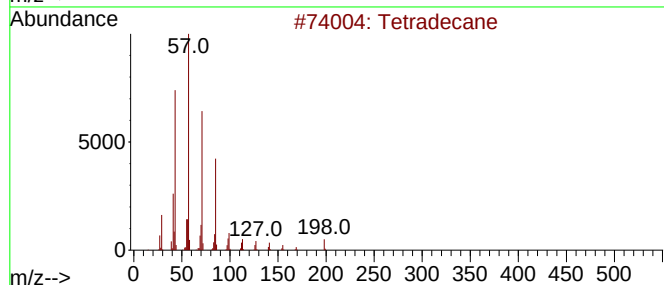
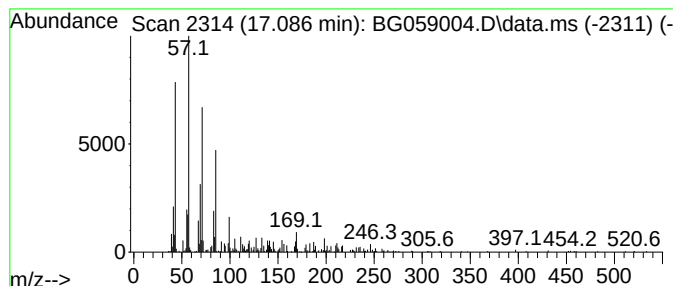
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.086	11.03 ng/ul	624851	Phenanthrene-d10		17.697	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tetradecane		198	C14H30	000629-59-4	81
2	Pentadecane		212	C15H32	000629-62-9	76
3	Hexadecane, 2,6,10,14-tetramethyl-		282	C20H42	000638-36-8	72
4	Octadecane, 2-methyl-		268	C19H40	001560-88-9	68
5	Tetradecane		198	C14H30	000629-59-4	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

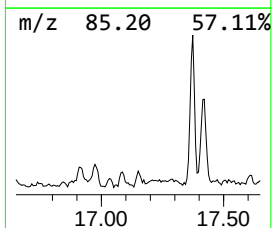
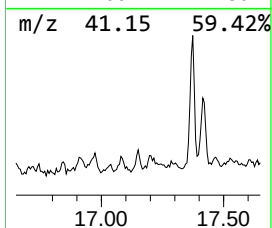
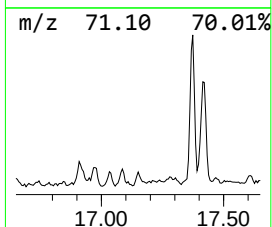
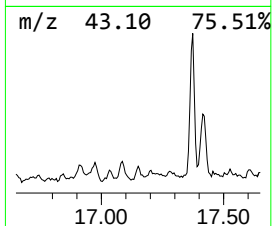
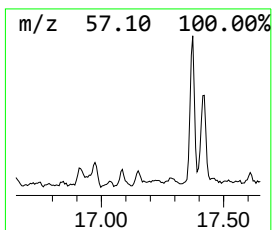
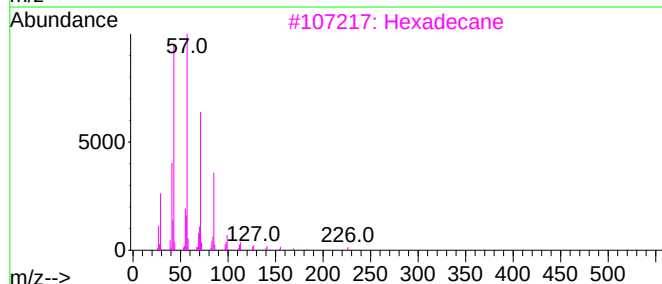
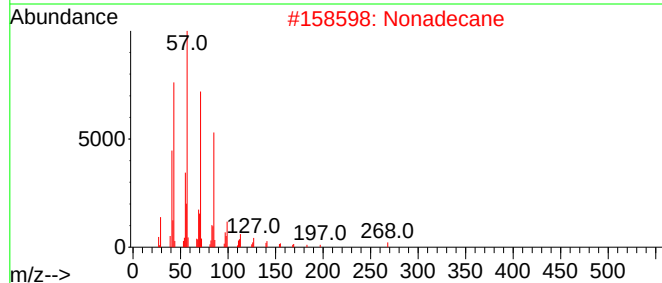
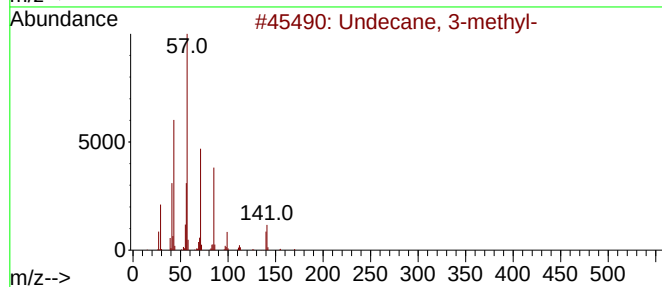
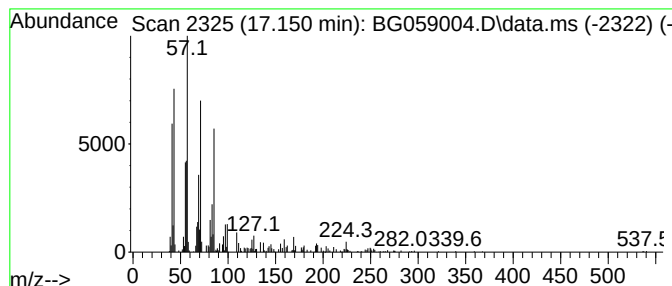
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.150	7.63 ng/ul	432290	Phenanthrene-d10	17.697

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 3-methyl-	170	C12H26	001002-43-3	74
2		Nonadecane	268	C19H40	000629-92-5	58
3		Hexadecane	226	C16H34	000544-76-3	53
4		Heptadecane, 3-methyl-	254	C18H38	006418-44-6	52
5		Octacosane	394	C28H58	000630-02-4	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

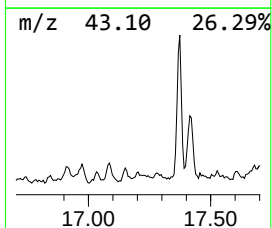
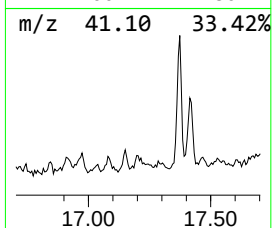
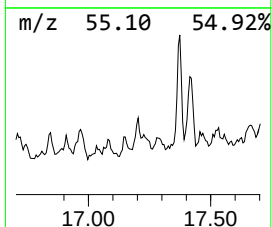
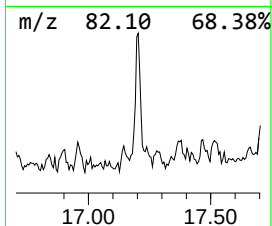
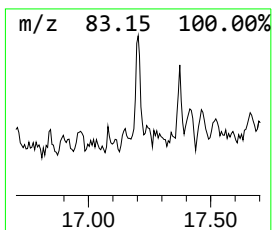
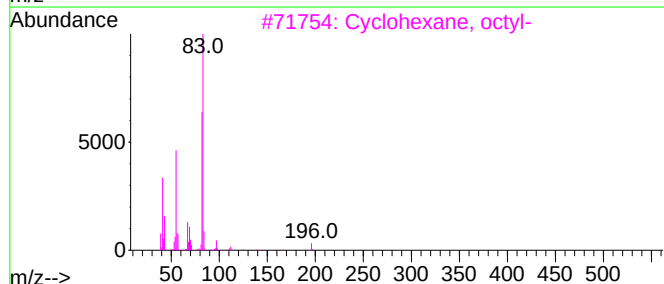
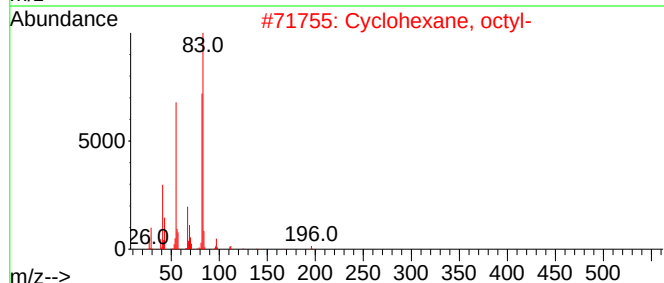
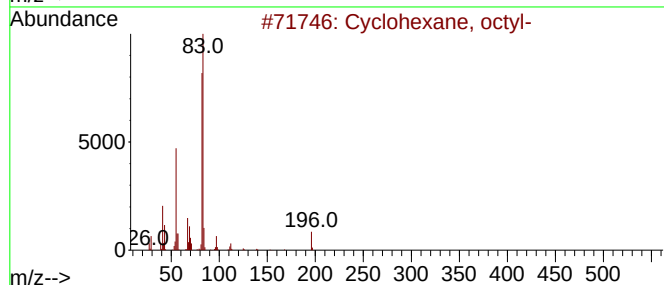
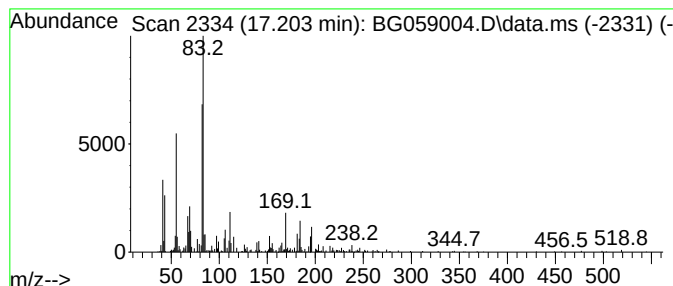
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 (DEL) Alkane: Cyclic17.203 Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.203	15.98 ng/ul	905155	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, octyl-	196	C14H28	001795-15-9	68
2			Cyclohexane, octyl-	196	C14H28	001795-15-9	68
3			Cyclohexane, octyl-	196	C14H28	001795-15-9	62
4			Cyclohexane, undecyl-	238	C17H34	054105-66-7	58
5			Cyclohexane, pentyl-	154	C11H22	004292-92-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

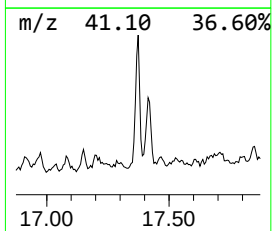
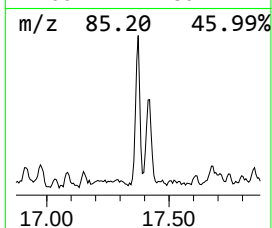
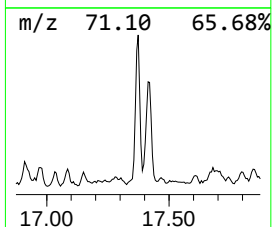
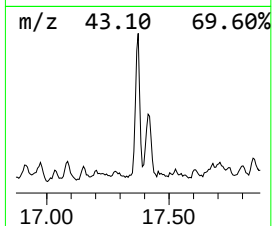
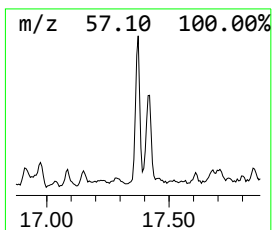
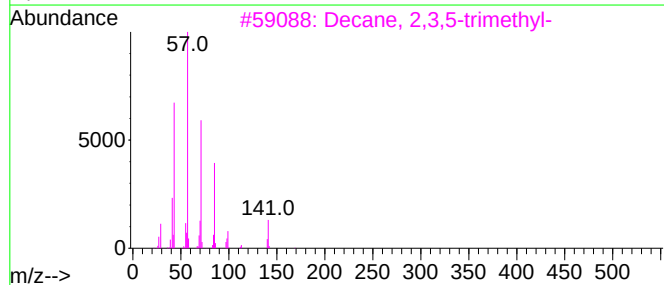
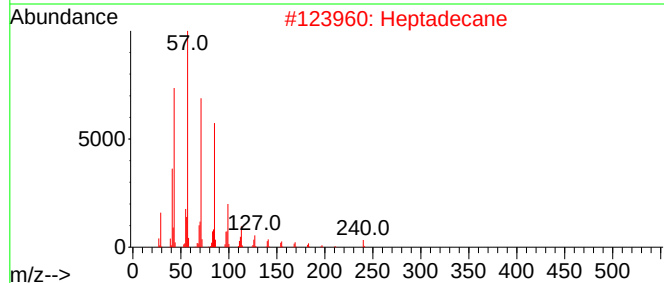
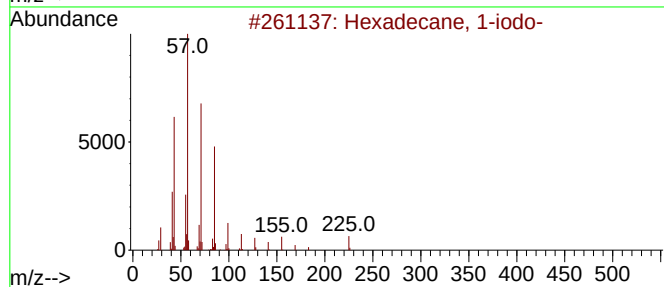
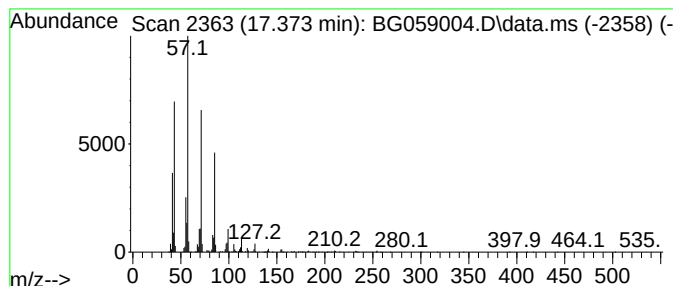
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 68 Hexadecane, 1-iodo- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.374	62.38 ng/ul	3533200	Phenanthrene-d10	17.697	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	90
2	Heptadecane	240	C17H36	000629-78-7	90
3	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
4	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
5	Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

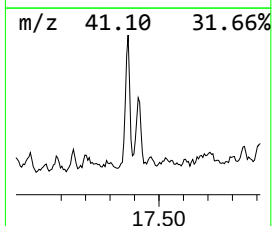
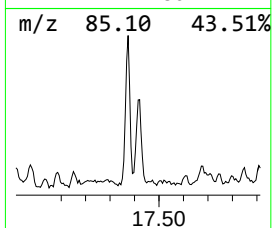
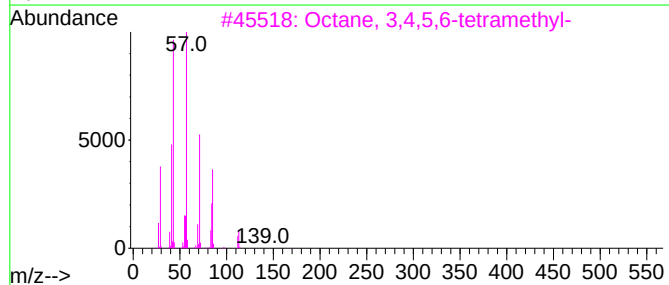
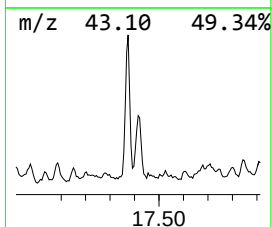
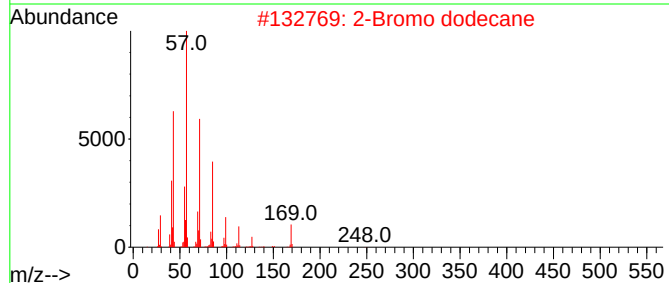
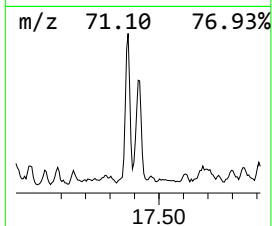
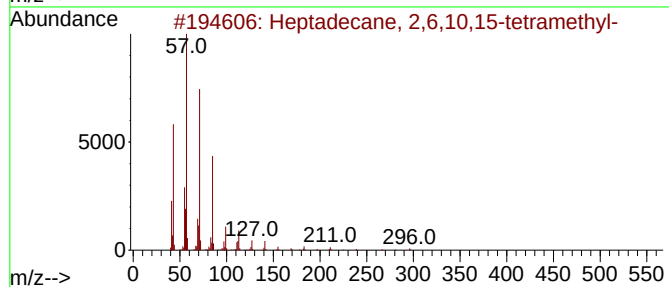
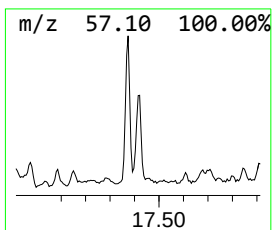
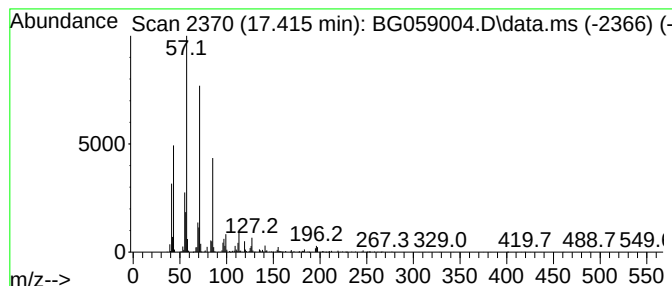
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 69 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.415	46.88 ng/ul	2655370	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	86
3			Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	72
4			Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	72
5			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

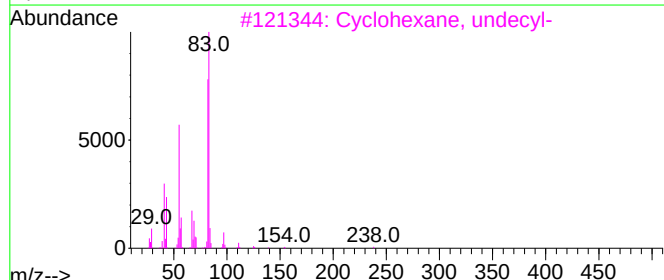
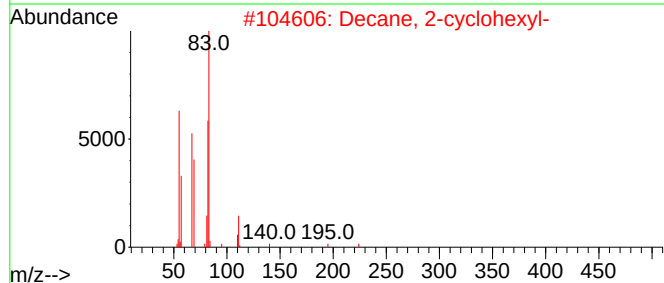
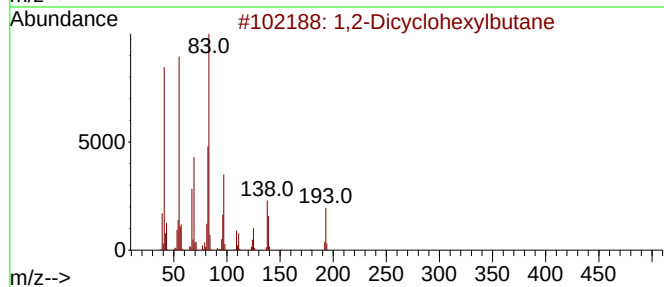
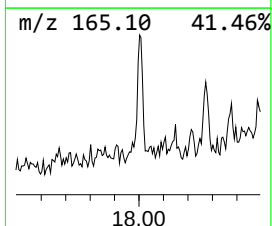
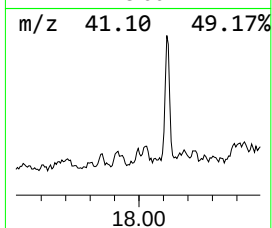
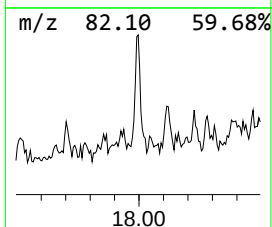
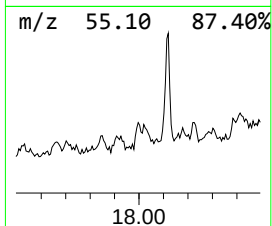
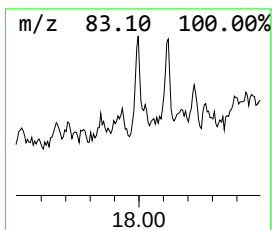
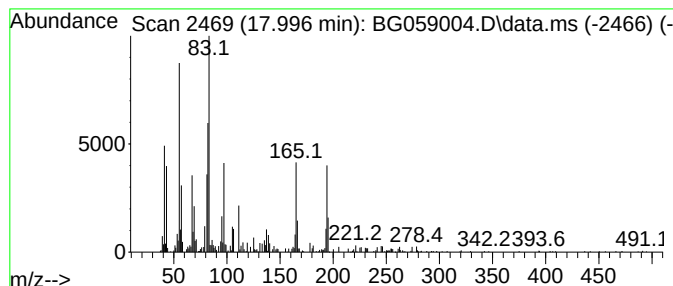
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 70 unknown-07 Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.996	8.43 ng/ul	477377	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Dicyclohexylbutane	222	C16H30	054890-01-6	43
2			Decane, 2-cyclohexyl-	224	C16H32	013151-73-0	38
3			Cyclohexane, undecyl-	238	C17H34	054105-66-7	30
4			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	27
5			3-Cyclohexyl-1-chloropropane	160	C9H17Cl	001124-62-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

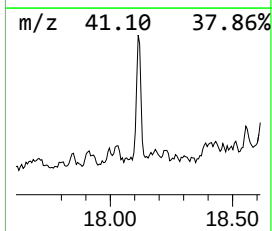
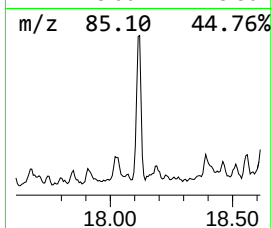
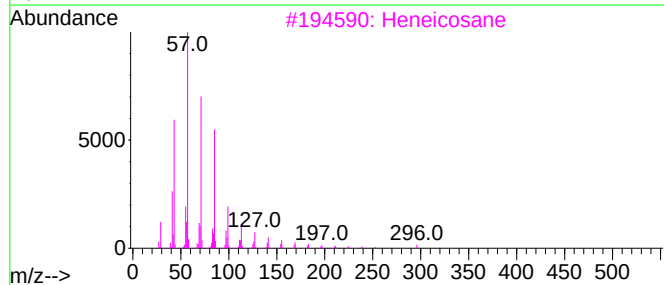
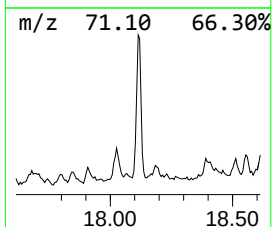
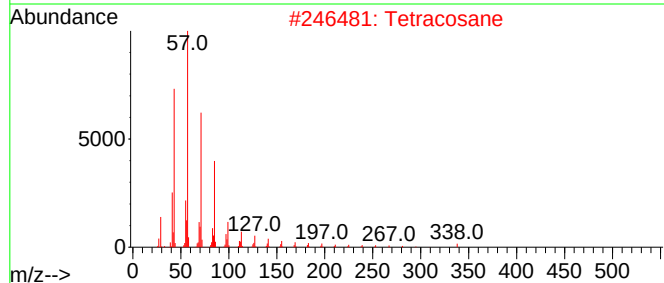
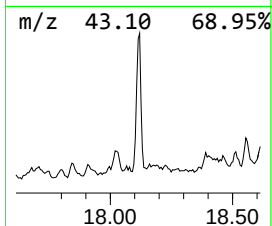
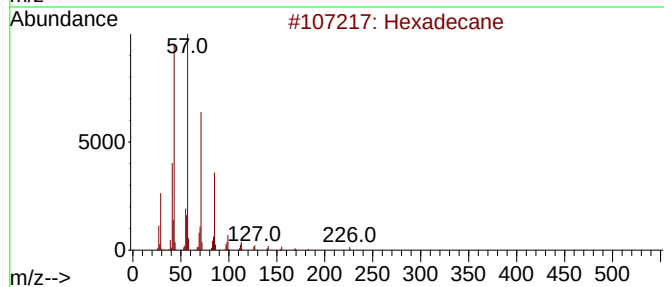
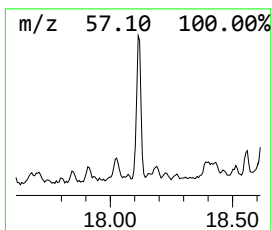
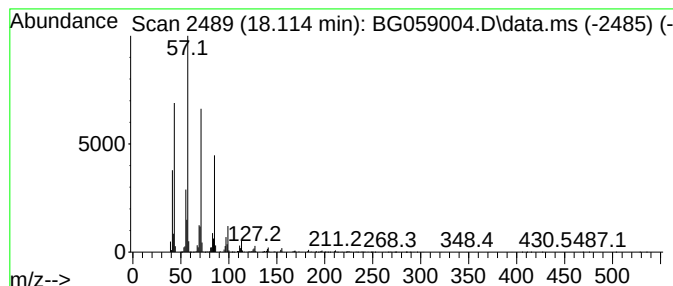
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 71 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.114	79.48 ng/ul	4502030	Phenanthrene-d10	17.697

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	90
2		Tetracosane	338	C24H50	000646-31-1	90
3		Heneicosane	296	C21H44	000629-94-7	90
4		Tetradecane	198	C14H30	000629-59-4	86
5		Hexacosane	366	C26H54	000630-01-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

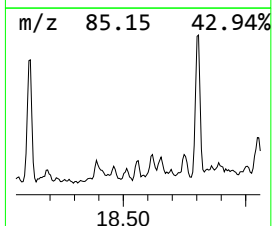
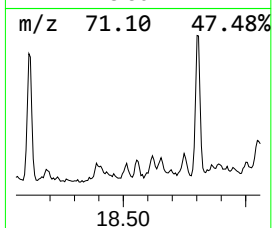
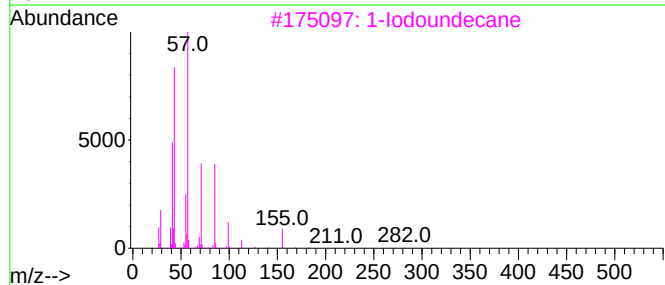
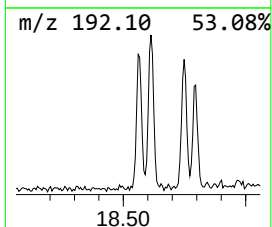
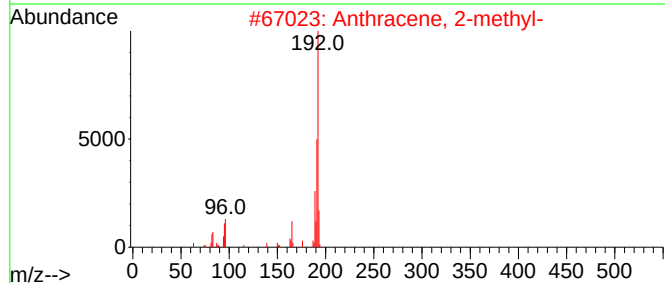
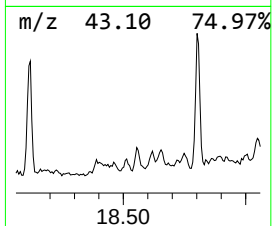
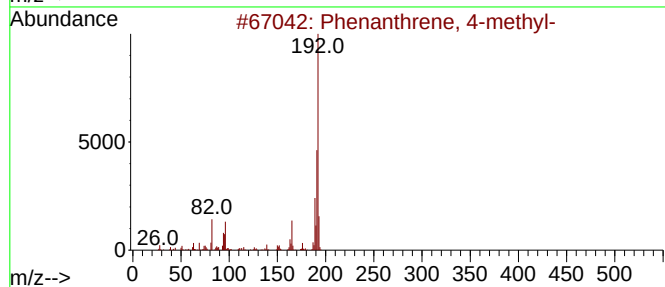
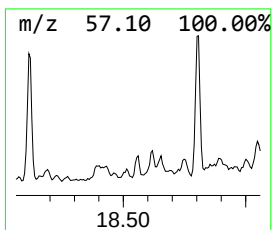
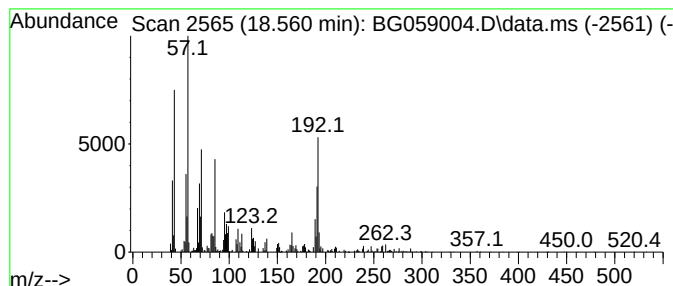
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 72 Phenanthrene, 4-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.560	25.31 ng/ul	1433780	Phenanthrene-d10	17.697	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	80
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	42
3	1-Iodoundecane	282	C11H23I	004282-44-4	42
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	42
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

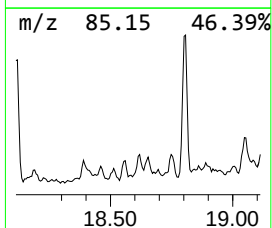
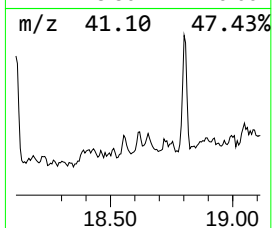
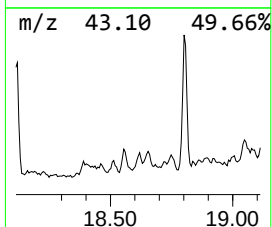
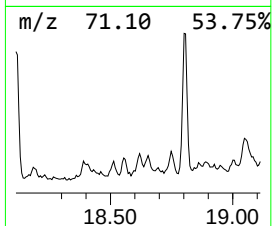
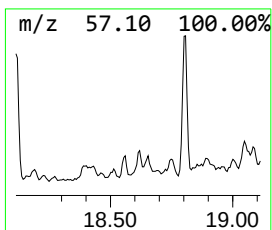
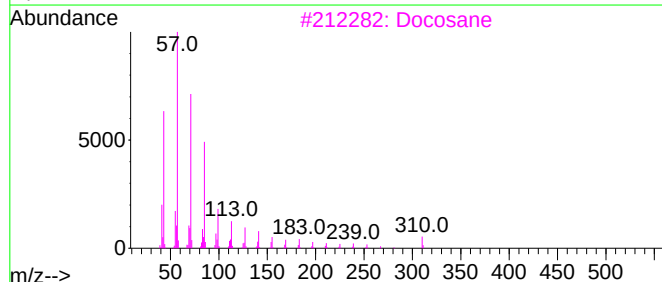
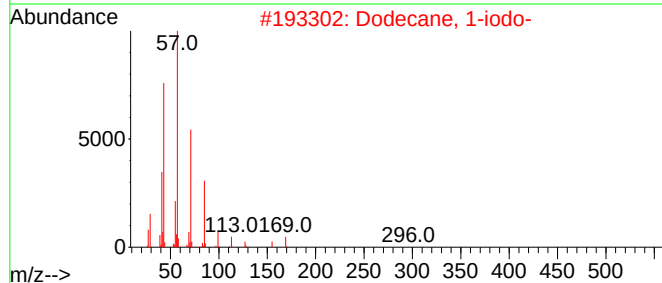
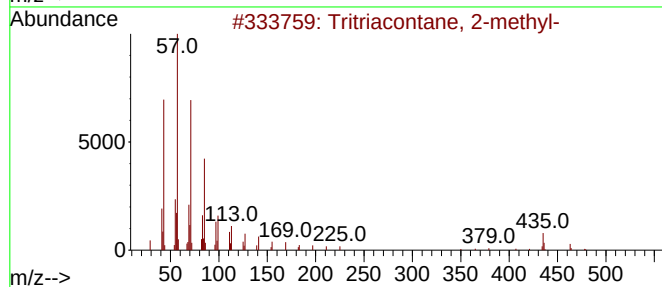
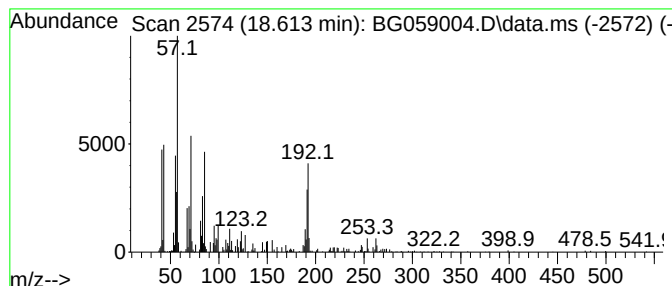
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 73 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.613	13.61 ng/ul	770941	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tritriacontane, 2-methyl-	479	C34H70	066214-27-5	64
2			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	41
3			Docosane	310	C22H46	000629-97-0	38
4			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	38
5			Octadecane	254	C18H38	000593-45-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

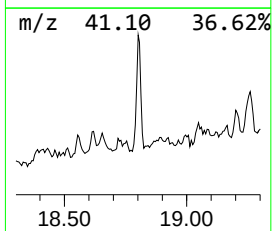
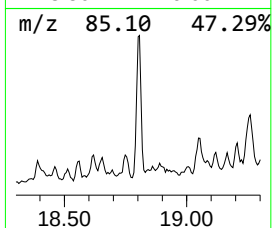
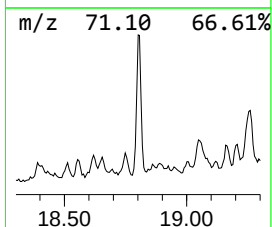
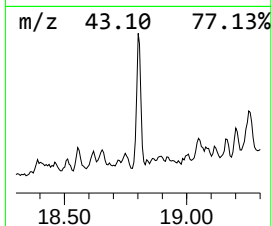
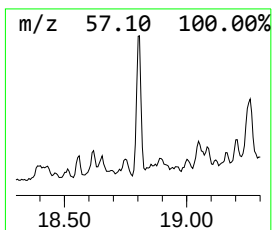
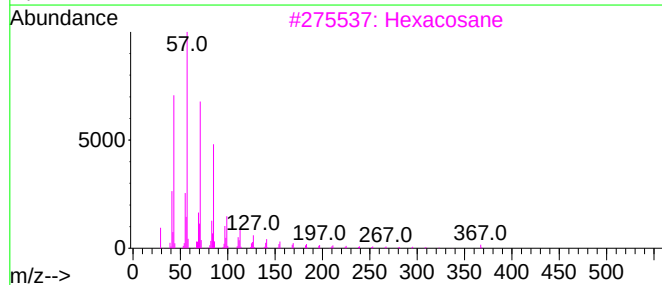
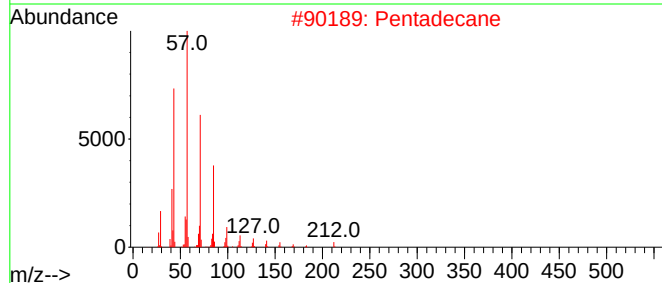
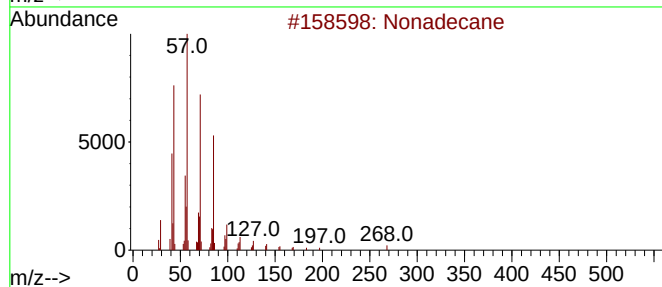
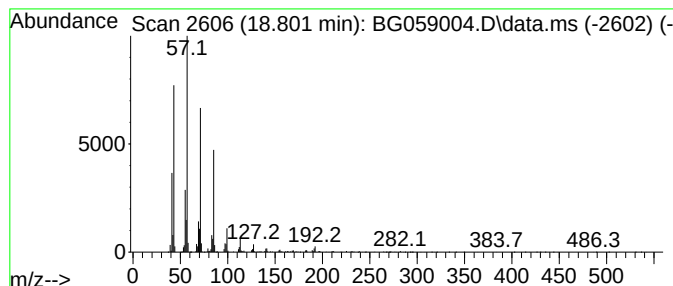
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 74 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.801	86.04 ng/ul	4873190	Phenanthrene-d10	17.697

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	93
2			Pentadecane	212	C15H32	000629-62-9	90
3			Hexacosane	366	C26H54	000630-01-3	90
4			Heptadecane	240	C17H36	000629-78-7	90
5			Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

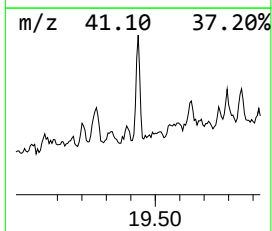
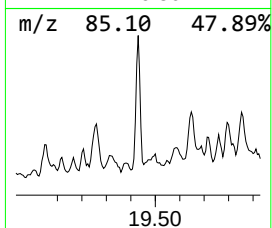
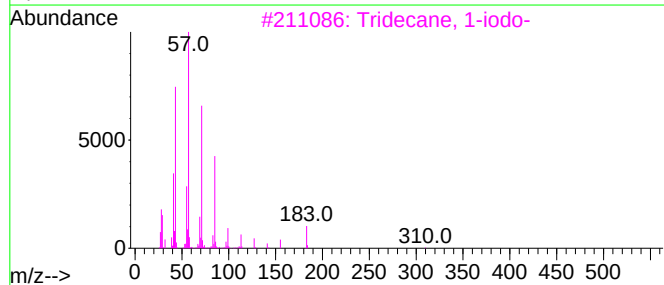
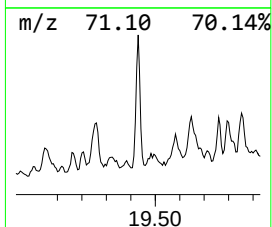
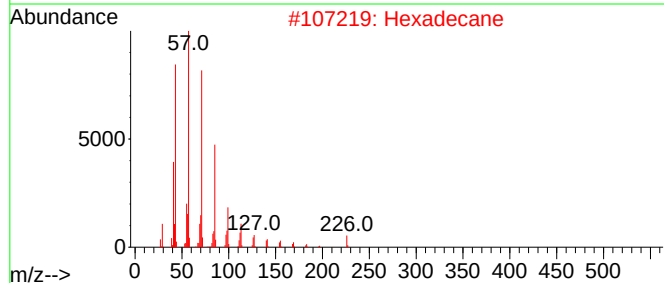
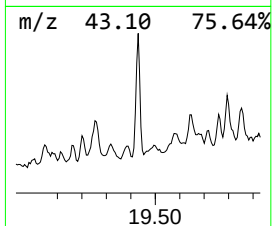
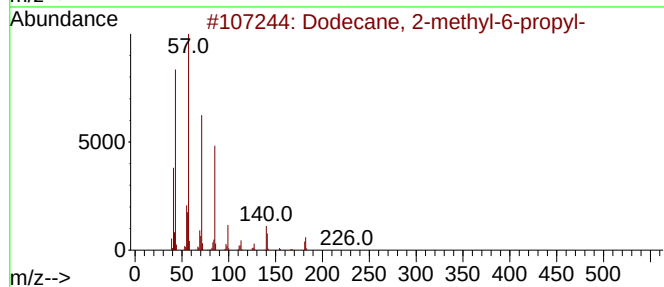
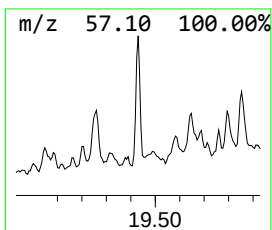
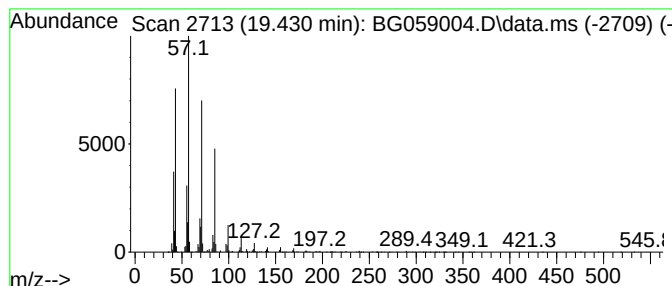
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 75 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.430	83.34 ng/ul	4720190	Phenanthrene-d10	17.697	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	94
2	Hexadecane	226	C16H34	000544-76-3	91
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
4	Tridecane, 7-propyl-	226	C16H34	055045-09-5	87
5	Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

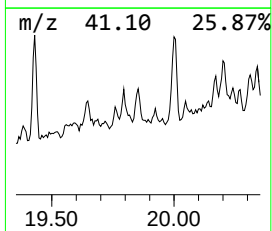
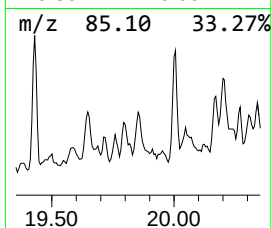
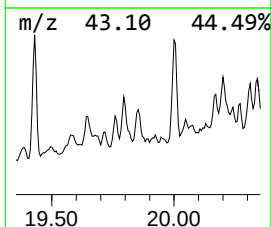
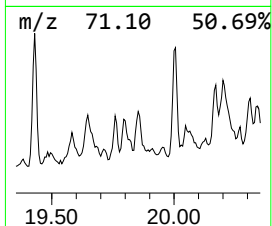
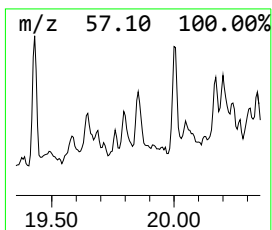
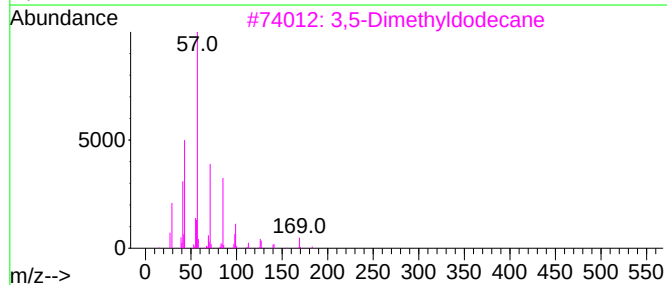
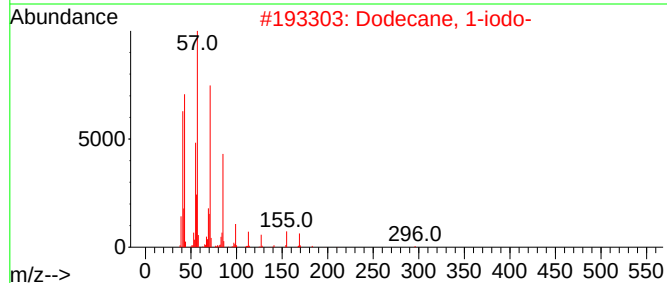
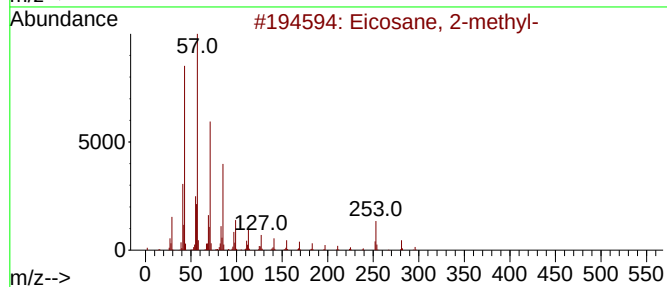
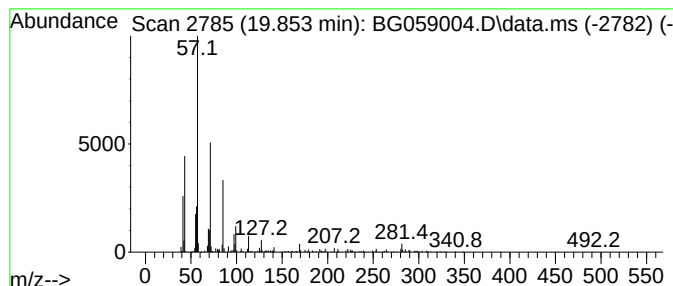
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 76 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.853	34.45 ng/ul	1951420	Phenanthrene-d10		17.697	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane, 2-methyl-		296	C21H44	001560-84-5	87
2	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	86
3	3,5-Dimethyldodecane		198	C14H30	107770-99-0	81
4	Tridecane, 2-methyl-		198	C14H30	001560-96-9	80
5	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
EZYJ9

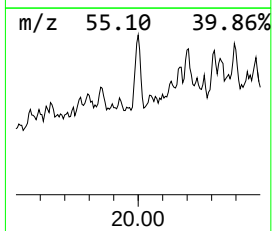
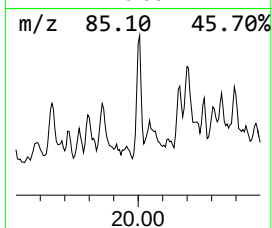
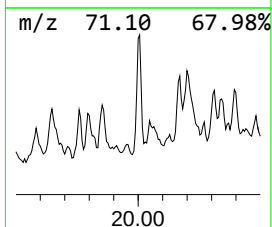
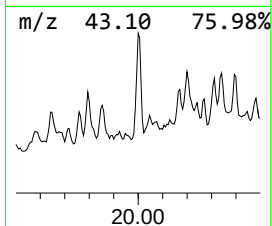
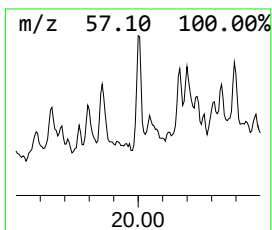
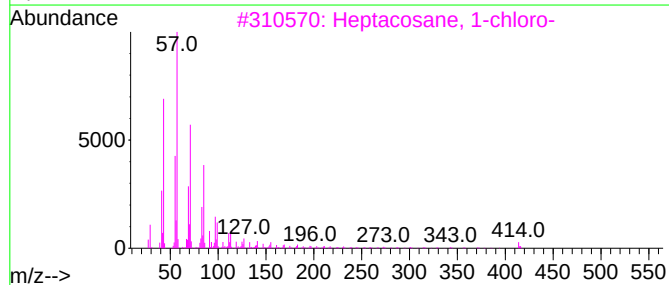
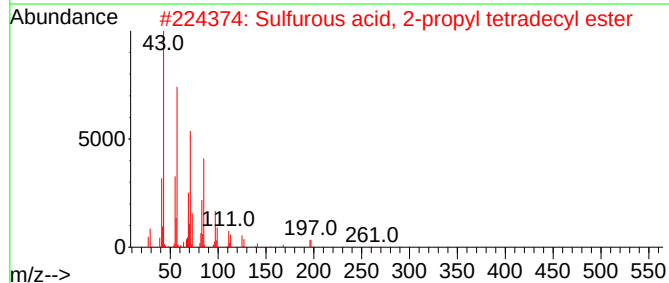
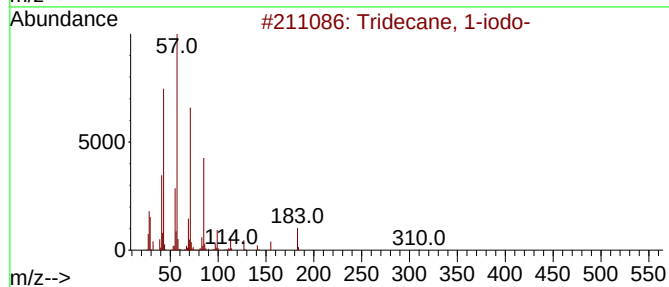
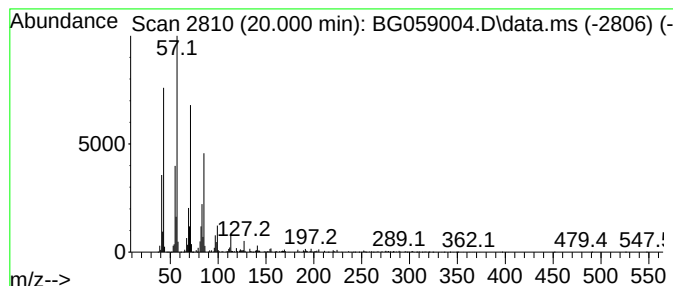
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 77 Tridecane, 1-iodo- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.000	16.38 ng/ul	5128280	Chrysene-d12	22.033

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
2			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	91
3			Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	90
4			Tetratetracontane	619	C44H90	007098-22-8	90
5			Heptacosane	380	C27H56	000593-49-7	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

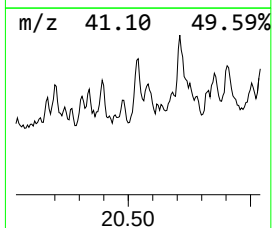
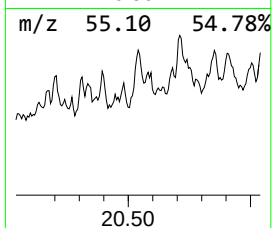
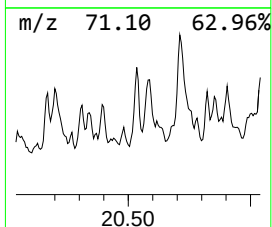
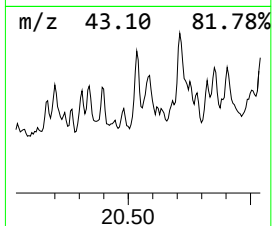
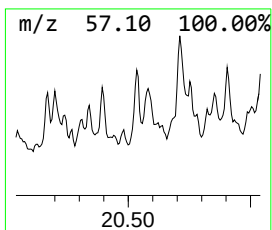
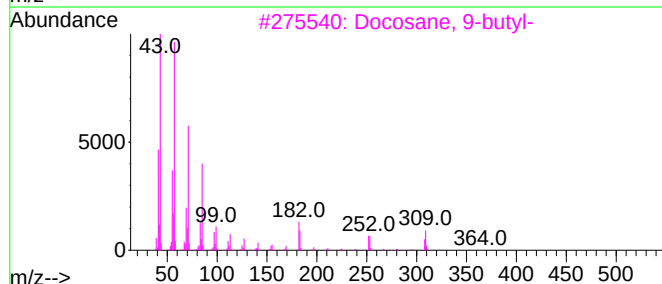
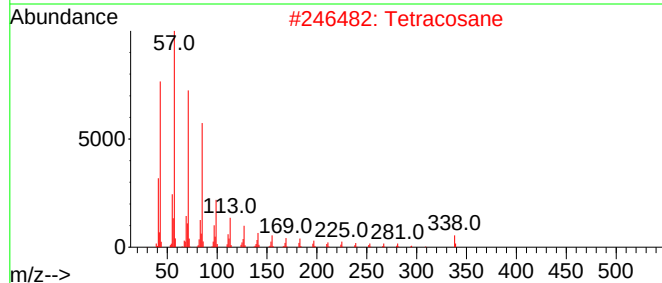
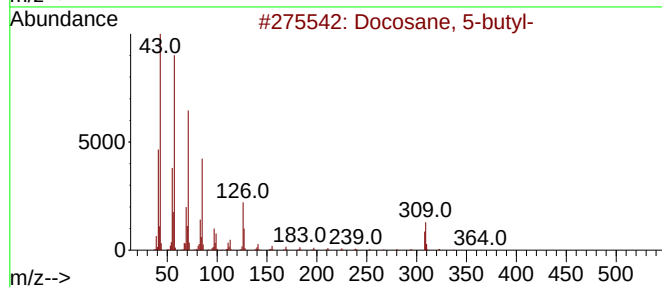
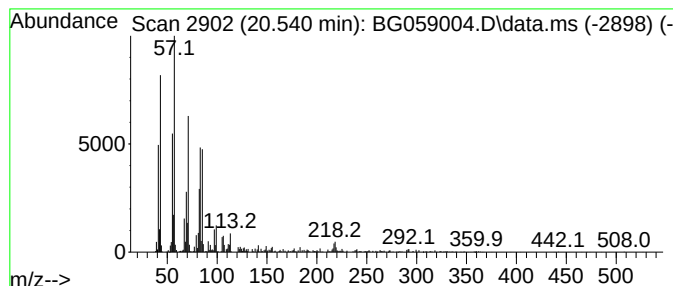
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 78 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.540	14.26 ng/ul	4463970	Chrysene-d12	22.033	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 5-butyl-	366	C26H54	055282-16-1	95
2	Tetracosane	338	C24H50	000646-31-1	78
3	Docosane, 9-butyl-	366	C26H54	055282-14-9	70
4	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	70
5	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

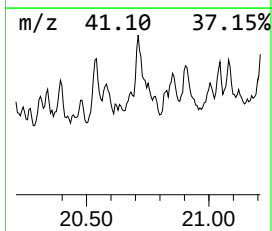
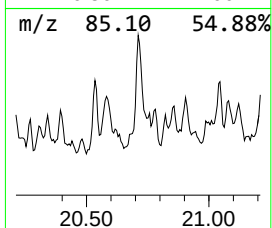
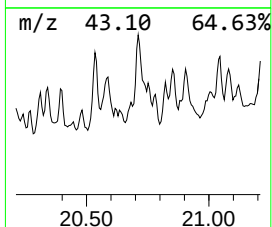
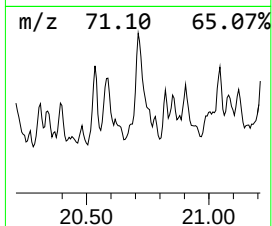
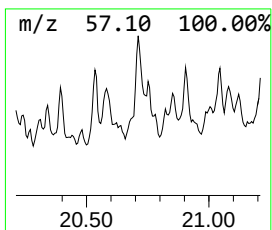
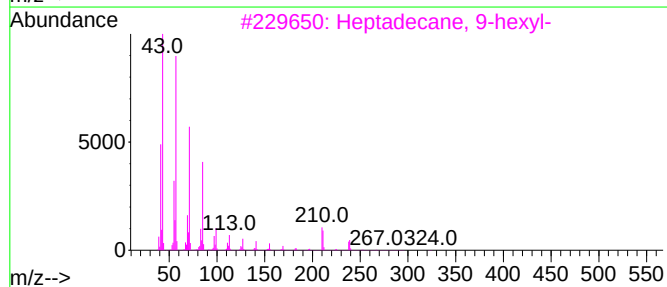
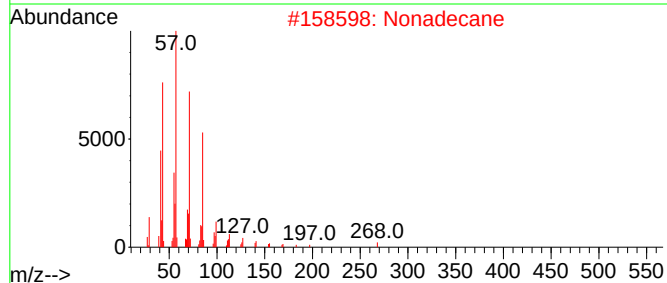
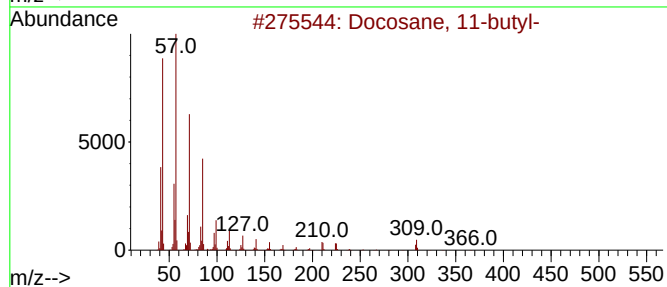
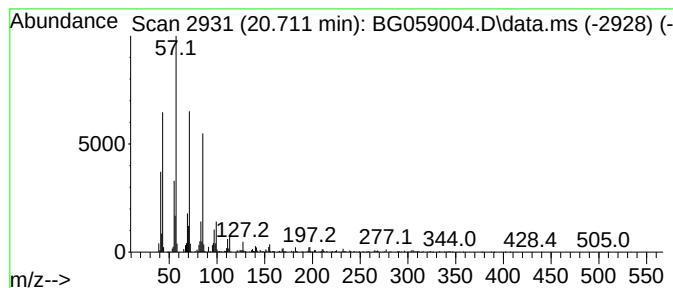
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 79 (DEL) Alkane: Straight-Chai... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD		R.T.	
20.711	16.53 ng/ul	5176960	Chrysene-d12		22.033	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-		366	C26H54	013475-76-8	93
2	Nonadecane		268	C19H40	000629-92-5	91
3	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	91
4	2-Bromo dodecane		248	C12H25Br	013187-99-0	91
5	Octadecane, 5,14-dibutyl-		366	C26H54	055282-13-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
Data File : BG059004.D
Acq On : 12 Sep 2023 6:29
Operator : MA/JU
Sample : 04235-08
Misc :
ALS Vial : 7 Sample Multiplier: 1

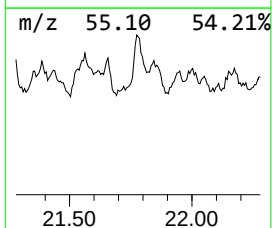
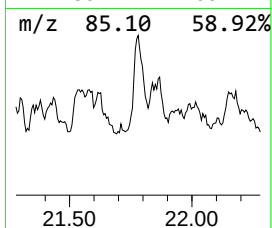
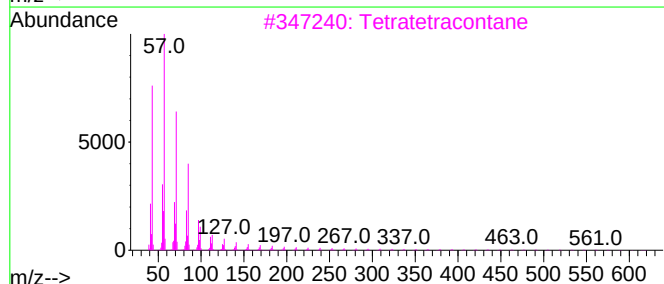
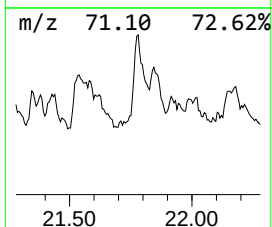
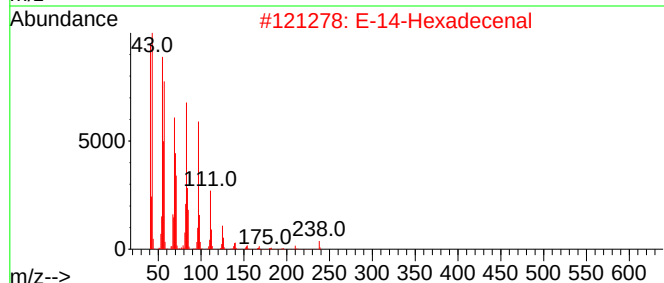
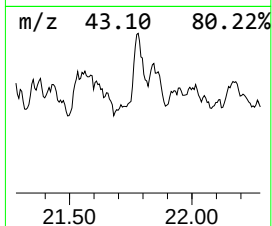
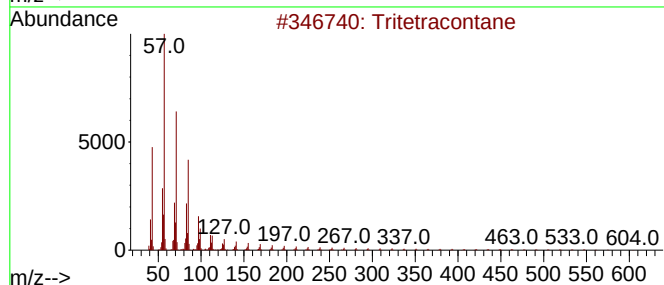
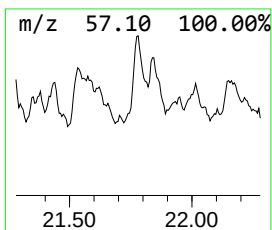
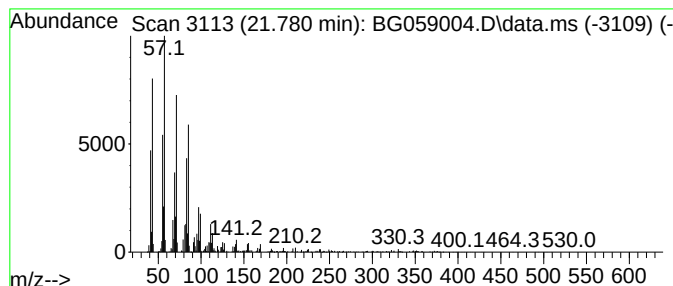
Instrument :
BNA_G
ClientSampleId :
EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 80 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.780	20.00 ng/ul	6262330	Chrysene-d12	22.033	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tritetracontane	605	C43H88	007098-21-7	83
2	E-14-Hexadecenal	238	C16H30O	330207-53-9	80
3	Tetratetracontane	619	C44H90	007098-22-8	72
4	Hexacosane	366	C26H54	000630-01-3	68
5	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG059004.D
 Acq On : 12 Sep 2023 6:29
 Operator : MA/JU
 Sample : 04235-08
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EZYJ9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	8.913	3.6	ng/uI	131864	1	8.325	733569	20.0
3-Fluoro-2-form...	9.383	8.3	ng/uI	306338	1	8.325	733569	20.0
(DEL) Alkane: S...	9.741	6.1	ng/uI	224834	1	8.325	733569	20.0
(DEL) Alkane: S...	9.900	3.0	ng/uI	264644	2	11.169	1745550	20.0
trans-Decalin, ...	10.035	3.9	ng/uI	336234	2	11.169	1745550	20.0
(DEL) Alkane: C...	10.211	5.1	ng/uI	448592	2	11.169	1745550	20.0
unknown-01	10.305	10.8	ng/uI	938141	2	11.169	1745550	20.0
(DEL) Alkane: S...	10.405	4.6	ng/uI	399142	2	11.169	1745550	20.0
(DEL) Alkane: S...	10.493	4.4	ng/uI	381820	2	11.169	1745550	20.0
(DEL) Alkane: S...	10.593	4.9	ng/uI	423750	2	11.169	1745550	20.0
unknown-02	10.799	3.6	ng/uI	310953	2	11.169	1745550	20.0
(DEL) Alkane: C...	10.940	4.0	ng/uI	346799	2	11.169	1745550	20.0
(DEL) Alkane: C...	10.993	5.0	ng/uI	438027	2	11.169	1745550	20.0
(DEL) Alkane: S...	11.046	42.8	ng/uI	3738180	2	11.169	1745550	20.0
unknown-03	11.093	3.9	ng/uI	341387	2	11.169	1745550	20.0
(DEL) Alkane: S...	11.228	28.0	ng/uI	2442620	2	11.169	1745550	20.0
(DEL) Alkane: C...	11.592	11.1	ng/uI	971202	2	11.169	1745550	20.0
(DEL) Alkane: C...	11.804	23.8	ng/uI	2080480	2	11.169	1745550	20.0
(DEL) Alkane: S...	11.909	7.7	ng/uI	669200	2	11.169	1745550	20.0
(DEL) Alkane: S...	11.986	10.8	ng/uI	939714	2	11.169	1745550	20.0
(DEL) Alkane: S...	12.086	38.9	ng/uI	3395880	2	11.169	1745550	20.0
Zinc, bis[2-(1,...	12.144	7.2	ng/uI	632199	2	11.169	1745550	20.0
unknown-04	12.303	4.4	ng/uI	380868	2	11.169	1745550	20.0
(DEL) Alkane: S...	12.362	9.2	ng/uI	800430	2	11.169	1745550	20.0
(DEL) Alkane: S...	12.426	8.8	ng/uI	766116	2	11.169	1745550	20.0
(DEL) Alkane: S...	12.485	111.0	ng/uI	9684610	2	11.169	1745550	20.0
unknown-05	12.556	7.6	ng/uI	661293	2	11.169	1745550	20.0
Sulfurous acid,...	12.608	3.6	ng/uI	314056	2	11.169	1745550	20.0
(DEL) Alkane: S...	12.691	21.8	ng/uI	1901780	2	11.169	1745550	20.0
(DEL) Alkane: C...	12.896	12.5	ng/uI	1094700	2	11.169	1745550	20.0
Ethyl octacosyl...	13.102	23.0	ng/uI	1218560	3	14.947	1061220	20.0
Pentane, 3-(bro...	13.137	15.5	ng/uI	821173	3	14.947	1061220	20.0
(DEL) Alkane: C...	13.190	48.5	ng/uI	2571370	3	14.947	1061220	20.0
Dodecane, 1-iodo-	13.278	34.9	ng/uI	1854170	3	14.947	1061220	20.0
(DEL) Alkane: C...	13.331	2.7	ng/uI	141071	3	14.947	1061220	20.0
(DEL) Alkane: S...	13.366	24.1	ng/uI	1278530	3	14.947	1061220	20.0
Pentacosane	13.413	83.4	ng/uI	4423480	3	14.947	1061220	20.0
(DEL) Alkane: S...	13.707	252.5	ng/uI	13395800	3	14.947	1061220	20.0
(DEL) Alkane: S...	13.860	6.5	ng/uI	343008	3	14.947	1061220	20.0
(DEL) Alkane: S...	13.942	9.0	ng/uI	476246	3	14.947	1061220	20.0
Naphthalene, 1,...	14.101	14.4	ng/uI	762976	3	14.947	1061220	20.0
(DEL) Alkane: S...	14.224	8.8	ng/uI	467479	3	14.947	1061220	20.0
Naphthalene, 1,...	14.265	38.1	ng/uI	2023800	3	14.947	1061220	20.0
(DEL) Alkane: S...	14.471	14.7	ng/uI	780141	3	14.947	1061220	20.0
(DEL) Alkane: S...	14.771	227.4	ng/uI	12067100	3	14.947	1061220	20.0
Sulfurous acid,...	14.853	18.3	ng/uI	970926	3	14.947	1061220	20.0
unknown-06	15.017	5.6	ng/uI	297383	3	14.947	1061220	20.0
(DEL) Alkane: S...	15.205	13.8	ng/uI	731422	3	14.947	1061220	20.0
(DEL) Alkane: S...	15.258	18.7	ng/uI	994079	3	14.947	1061220	20.0
4,6,8-Trimethyl...	15.317	22.4	ng/uI	1190380	3	14.947	1061220	20.0
(DEL) Alkane: S...	15.376	39.3	ng/uI	2086830	3	14.947	1061220	20.0
(DEL) Alkane: S...	15.446	15.5	ng/uI	821256	3	14.947	1061220	20.0

Naphthalene, 2,...	15.540	10.3	ng/ul	545971	3	14.947	1061220	20.0
Naphthalene, 1,...	15.599	25.0	ng/ul	1328830	3	14.947	1061220	20.0
(DEL) Alkane: S...	15.717	169.1	ng/ul	8971520	3	14.947	1061220	20.0
Octadecane, 1-c...	15.770	11.8	ng/ul	626493	3	14.947	1061220	20.0
(DEL) Alkane: C...	15.828	6.5	ng/ul	347271	3	14.947	1061220	20.0
Naphthalene, 2-...	16.016	10.8	ng/ul	571881	3	14.947	1061220	20.0
(DEL) Alkane: S...	16.110	74.8	ng/ul	3966650	3	14.947	1061220	20.0
(DEL) Alkane: S...	16.157	4.8	ng/ul	254208	3	14.947	1061220	20.0
(DEL) Alkane: S...	16.269	11.8	ng/ul	623905	3	14.947	1061220	20.0
(DEL) Alkane: C...	16.339	17.1	ng/ul	969264	4	17.697	1132800	20.0
(DEL) Alkane: S...	16.580	173.2	ng/ul	9807630	4	17.697	1132800	20.0
1-Methyl-3-pipe...	16.845	4.9	ng/ul	279438	4	17.697	1132800	20.0
(DEL) Alkane: S...	17.086	11.0	ng/ul	624851	4	17.697	1132800	20.0
(DEL) Alkane: S...	17.150	7.6	ng/ul	432290	4	17.697	1132800	20.0
(DEL) Alkane: C...	17.203	16.0	ng/ul	905155	4	17.697	1132800	20.0
Hexadecane, 1-i...	17.374	62.4	ng/ul	3533200	4	17.697	1132800	20.0
(DEL) Alkane: S...	17.415	46.9	ng/ul	2655370	4	17.697	1132800	20.0
unknown-07	17.996	8.4	ng/ul	477377	4	17.697	1132800	20.0
(DEL) Alkane: S...	18.114	79.5	ng/ul	4502030	4	17.697	1132800	20.0
Phenanthrene, 4...	18.560	25.3	ng/ul	1433780	4	17.697	1132800	20.0
(DEL) Alkane: S...	18.613	13.6	ng/ul	770941	4	17.697	1132800	20.0
(DEL) Alkane: S...	18.801	86.0	ng/ul	4873190	4	17.697	1132800	20.0
(DEL) Alkane: S...	19.430	83.3	ng/ul	4720190	4	17.697	1132800	20.0
(DEL) Alkane: S...	19.853	34.5	ng/ul	1951420	4	17.697	1132800	20.0
Tridecane, 1-iodo-	20.000	16.4	ng/ul	5128280	5	22.033	6262330	20.0
(DEL) Alkane: S...	20.540	14.3	ng/ul	4463970	5	22.033	6262330	20.0
(DEL) Alkane: S...	20.711	16.5	ng/ul	5176960	5	22.033	6262330	20.0
(DEL) Alkane: S...	21.780	20.0	ng/ul	6262330	5	22.033	6262330	20.0

Instrument :

BNA_G

ClientSampleId :

EZYJ9