

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062906.D
 Acq On : 18 Sep 2024 13:11
 Operator : JU/RC
 Sample : P3878-07ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDCJ0ME

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-BG091724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG062906.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.740	106	111	118	rBV	103605	188452	3.31%	0.521%
2	5.885	471	476	484	rVB	114016	182289	3.20%	0.504%
3	6.425	563	568	573	rBV3	39598	60772	1.07%	0.168%
4	6.860	638	642	646	rVV2	61265	94879	1.66%	0.262%
5	6.913	646	651	655	rVV2	71235	113592	1.99%	0.314%
6	6.954	655	658	662	rVV3	48046	70795	1.24%	0.196%
7	7.024	662	670	677	rVV2	274100	516425	9.06%	1.429%
8	7.189	692	698	703	rBV	135673	226987	3.98%	0.628%
9	7.389	723	732	743	rVV3	192075	409437	7.18%	1.133%
10	7.489	743	749	760	rVB2	343664	632780	11.10%	1.750%
11	7.777	791	798	803	rBV6	23844	59254	1.04%	0.164%
12	7.853	804	811	817	rVV2	224315	429077	7.53%	1.187%
13	7.918	817	822	828	rVB	161754	251757	4.42%	0.696%
14	7.976	828	832	839	rVB4	34519	58290	1.02%	0.161%
15	8.288	879	885	889	rBV4	32880	64128	1.12%	0.177%
16	8.393	898	903	909	rVB6	60371	122533	2.15%	0.339%
17	8.446	909	912	916	rVB2	45316	56836	1.00%	0.157%
18	8.517	916	924	927	rBV4	86949	213183	3.74%	0.590%
19	8.558	927	931	937	rVB	238998	383649	6.73%	1.061%
20	8.623	938	942	945	rBV	69148	93367	1.64%	0.258%
21	8.805	969	973	978	rBV3	35689	54998	0.96%	0.152%
22	8.858	978	982	989	rVB2	41548	72354	1.27%	0.200%
23	8.963	989	1000	1003	rBV3	65207	151666	2.66%	0.420%
24	9.004	1003	1007	1013	rVV	179906	312813	5.49%	0.865%
25	9.087	1015	1021	1026	rVB	407261	624369	10.95%	1.727%
26	9.204	1032	1041	1048	rBV10	56548	150250	2.64%	0.416%
27	9.498	1085	1091	1094	rBV6	26963	56638	0.99%	0.157%
28	9.645	1110	1116	1122	rVB3	48094	86228	1.51%	0.239%
29	9.727	1123	1130	1136	rBV3	186127	358828	6.29%	0.993%
30	9.792	1136	1141	1145	rBV6	68443	129929	2.28%	0.359%
31	9.933	1162	1165	1169	rVB2	66310	96887	1.70%	0.268%
32	10.003	1172	1177	1181	rBV5	38890	57964	1.02%	0.160%
33	10.185	1204	1208	1212	rBV3	36059	55048	0.97%	0.152%
34	10.268	1214	1222	1229	rVB	247408	461290	8.09%	1.276%
35	10.344	1229	1235	1239	rBV3	58960	96284	1.69%	0.266%
36	10.514	1259	1264	1268	rBV6	27723	59658	1.05%	0.165%

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Instrument :
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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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-BG091724.MA.M
 Title : SVOA CALIBRATION

37	10.632	1276	1284	1292	rBV2	418269	1049525	18.41%	2.903%
38	10.691	1292	1294	1301	rVB3	55782	93796	1.65%	0.259%
39	10.773	1301	1308	1320	rBV4	289400	682542	11.97%	1.888%
40	11.026	1345	1351	1357	rBV	295140	468724	8.22%	1.297%
41	11.149	1364	1372	1379	rBV5	63879	150568	2.64%	0.417%
42	11.560	1436	1442	1448	rVV10	40121	86572	1.52%	0.239%
43	11.678	1456	1462	1466	rBV2	125244	209570	3.68%	0.580%
44	11.742	1466	1473	1480	rVB3	190017	384398	6.74%	1.063%
45	11.907	1497	1501	1510	rVB	147151	242336	4.25%	0.670%
46	12.071	1523	1529	1532	rBV	239334	374688	6.57%	1.036%
47	12.107	1532	1535	1541	rVB	178136	241658	4.24%	0.668%
48	12.318	1564	1571	1577	rVB2	307374	539161	9.46%	1.491%
49	12.477	1594	1598	1603	rVV5	43531	81058	1.42%	0.224%
50	12.524	1603	1606	1616	rVB4	56208	119102	2.09%	0.329%
51	12.718	1634	1639	1642	rBV3	42980	74923	1.31%	0.207%
52	12.888	1664	1668	1677	rVB7	47874	91168	1.60%	0.252%
53	13.029	1688	1692	1699	rVB6	77463	119832	2.10%	0.331%
54	13.317	1736	1741	1748	rBV2	266820	394087	6.91%	1.090%
55	13.752	1811	1815	1819	rVB3	43058	60300	1.06%	0.167%
56	13.811	1821	1825	1827	rBV4	54242	90685	1.59%	0.251%
57	13.893	1835	1839	1845	rVB	464053	678660	11.90%	1.877%
58	13.975	1846	1853	1857	rBV3	99548	167266	2.93%	0.463%
59	14.122	1874	1878	1882	rBV2	113834	153474	2.69%	0.425%
60	14.181	1883	1888	1903	rVB	471370	814218	14.28%	2.252%
61	14.392	1920	1924	1929	rVB	608303	789824	13.85%	2.185%
62	14.486	1935	1940	1947	rVV	510374	795871	13.96%	2.202%
63	14.563	1950	1953	1959	rVV6	94631	167492	2.94%	0.463%
64	14.633	1960	1965	1970	rVV	334587	476254	8.35%	1.317%
65	14.692	1971	1975	1984	rVB3	232528	423460	7.43%	1.171%
66	15.115	2043	2047	2055	rVB	286285	468642	8.22%	1.296%
67	15.180	2056	2058	2062	rVB2	96840	110283	1.93%	0.305%
68	15.250	2066	2070	2078	rBV	322500	481861	8.45%	1.333%
69	15.344	2083	2086	2091	rVB	276518	365899	6.42%	1.012%
70	15.485	2105	2110	2115	rVB	624603	934212	16.39%	2.584%
71	15.597	2124	2129	2137	rVB2	236986	391292	6.86%	1.082%
72	15.750	2153	2155	2165	rVB	108707	169514	2.97%	0.469%
73	15.849	2167	2172	2177	rVB	145866	218442	3.83%	0.604%
74	16.120	2214	2218	2223	rVB3	143687	186783	3.28%	0.517%
75	16.208	2229	2233	2236	rBV	291046	416815	7.31%	1.153%
76	16.419	2267	2269	2280	rVB4	72598	118598	2.08%	0.328%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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77	16.554	2288	2292	2295	rBV4	72458	92430	1.62%	0.256%
78	16.895	2343	2350	2359	rBV	3528355	5701154	100.00%	15.771%
79	17.001	2365	2368	2373	rVB	229170	268820	4.72%	0.744%
80	17.054	2374	2377	2381	rVB3	118456	153407	2.69%	0.424%
81	17.242	2403	2409	2416	rBV	657308	1031897	18.10%	2.854%
82	17.336	2420	2425	2430	rVV2	705570	1019149	17.88%	2.819%
83	17.530	2454	2458	2465	rVB4	108000	204535	3.59%	0.566%
84	17.741	2490	2494	2499	rVB	215280	281154	4.93%	0.778%
85	17.853	2511	2513	2519	rVB7	70728	87298	1.53%	0.241%
86	17.912	2520	2523	2527	rVB2	64624	86469	1.52%	0.239%
87	18.135	2558	2561	2565	rBV5	42594	58872	1.03%	0.163%
88	18.435	2608	2612	2617	rBV	157344	187730	3.29%	0.519%
89	19.069	2716	2720	2722	rBV	114168	167948	2.95%	0.465%
90	19.175	2735	2738	2743	rBV6	49084	72799	1.28%	0.201%
91	19.633	2810	2816	2822	rBV	934986	1375952	24.13%	3.806%
92	20.180	2906	2909	2917	rVB5	76129	137404	2.41%	0.380%
93	20.673	2990	2993	2998	rVB2	68812	78999	1.39%	0.219%
94	21.161	3071	3076	3085	rVB3	230111	360214	6.32%	0.996%
95	21.490	3126	3132	3139	rBV	666146	1079661	18.94%	2.987%
96	21.684	3162	3165	3170	rVB4	47498	66921	1.17%	0.185%
97	22.260	3259	3263	3273	rVB4	76492	155774	2.73%	0.431%
98	22.912	3370	3374	3382	rVB4	37942	66482	1.17%	0.184%
99	24.316	3604	3613	3626	rVB	486905	1273246	22.33%	3.522%
100	24.527	3640	3649	3662	rVB	438150	1204854	21.13%	3.333%

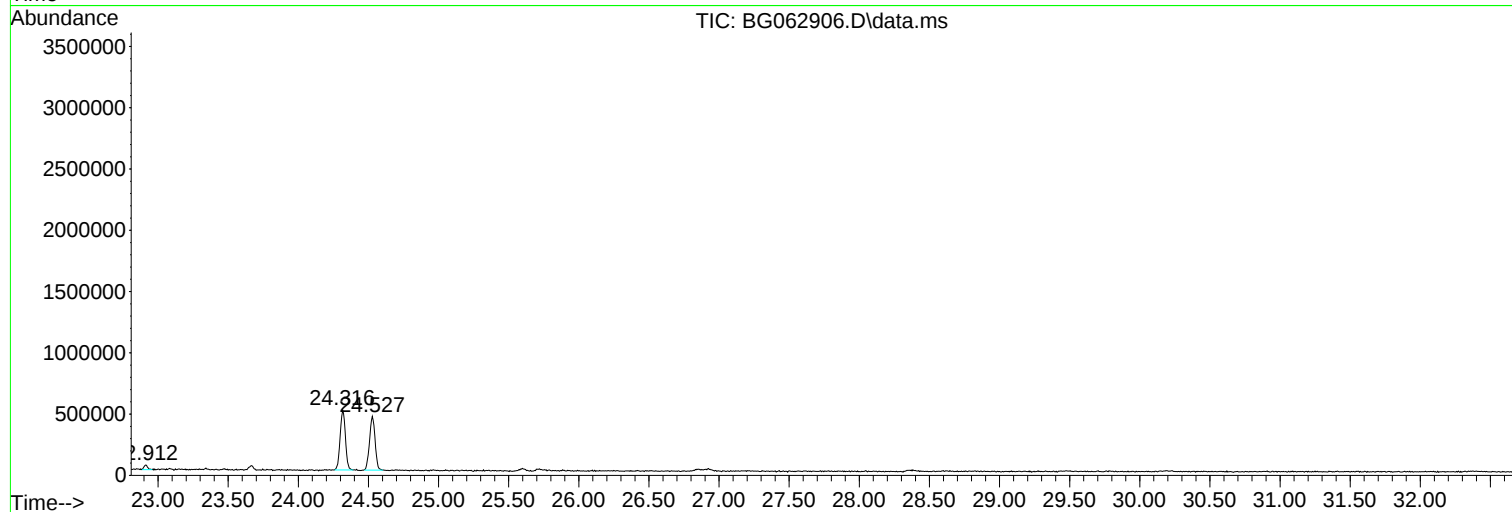
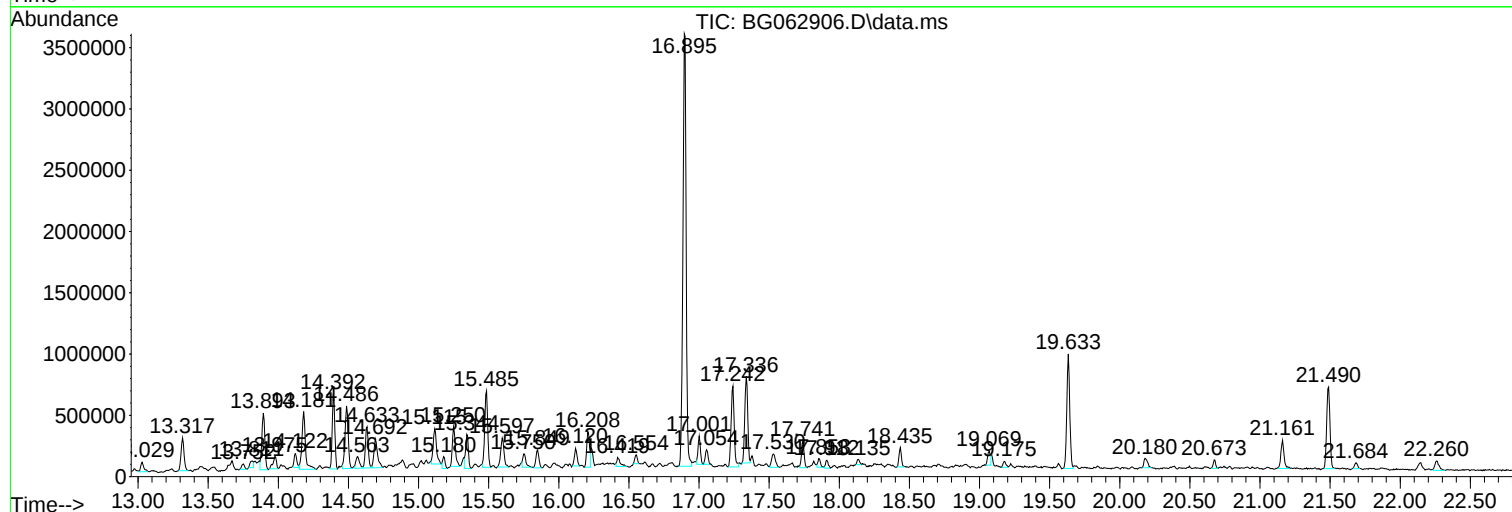
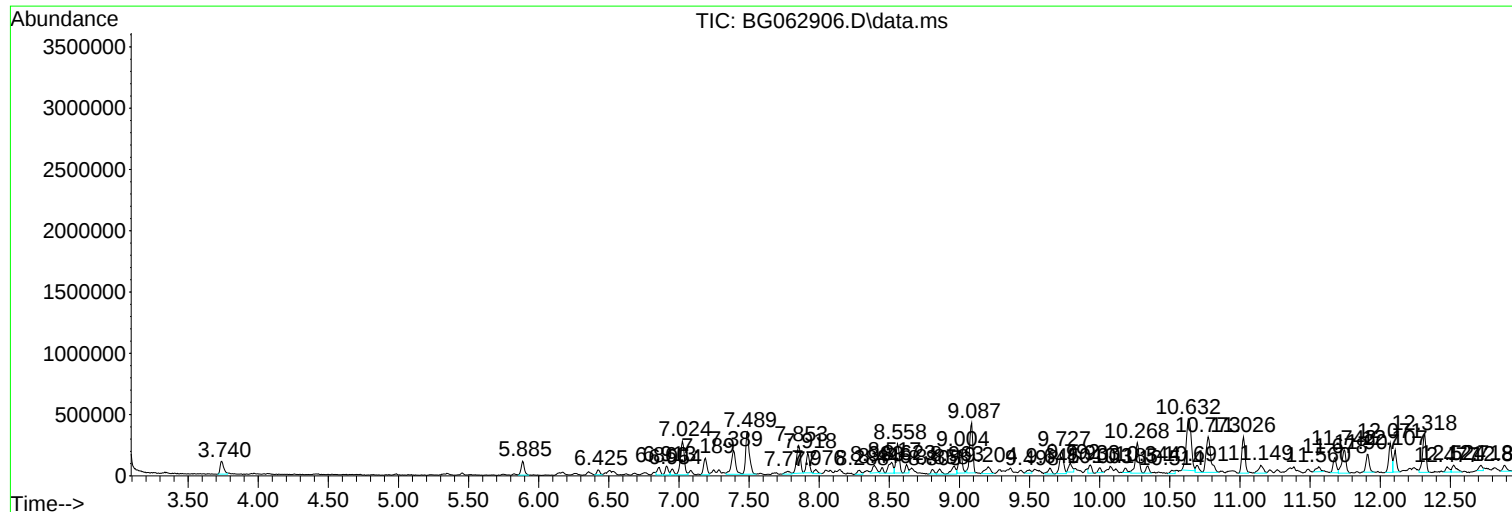
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Misc :
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Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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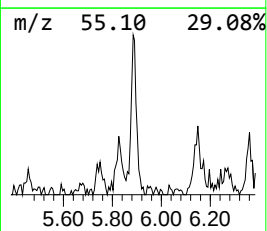
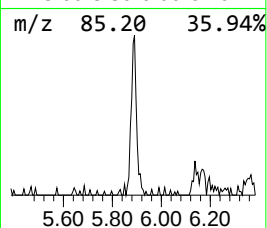
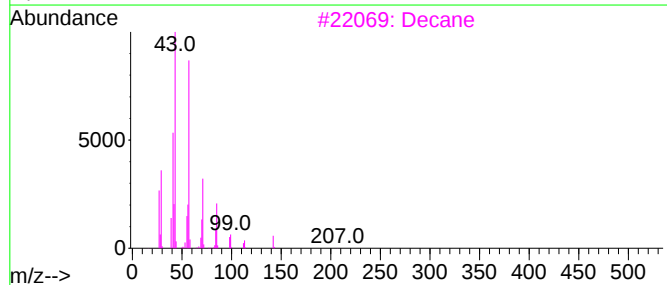
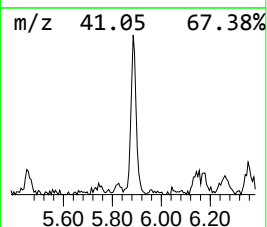
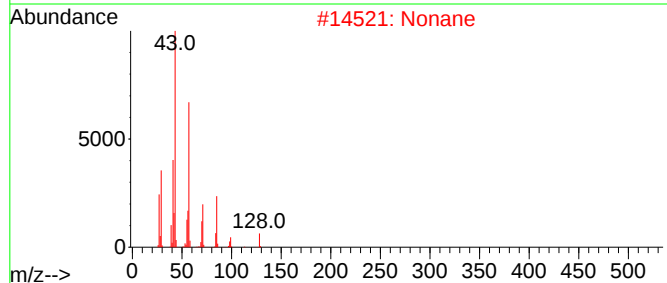
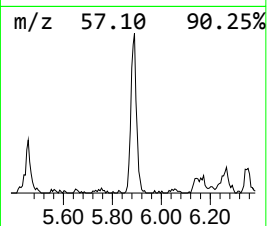
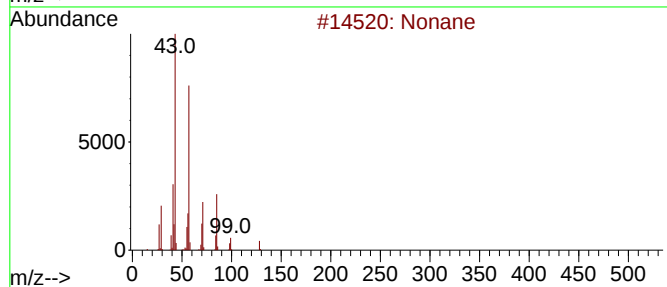
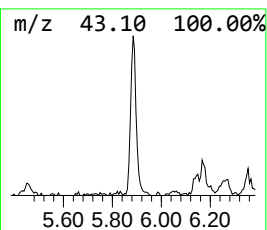
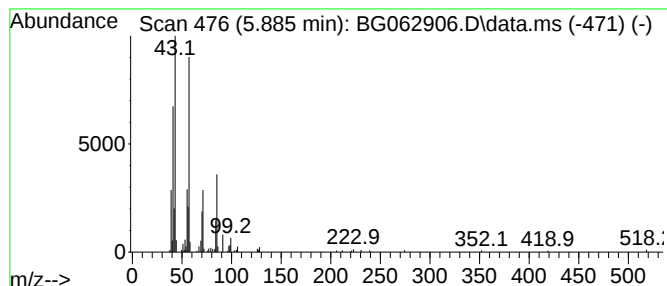
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.885	8.50 ng/ul	182289	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	93
2			Nonane	128	C9H20	000111-84-2	90
3			Decane	142	C10H22	000124-18-5	72
4			Decane	142	C10H22	000124-18-5	72
5			Nonane	128	C9H20	000111-84-2	64



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ClientSampleId :
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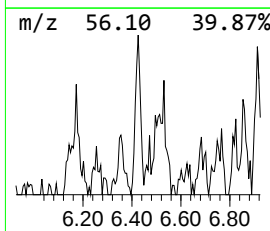
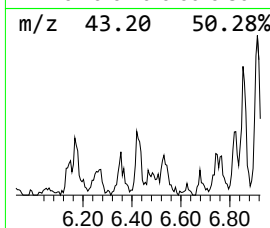
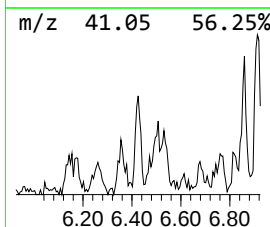
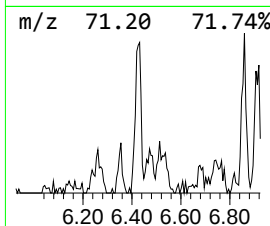
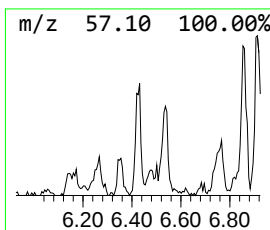
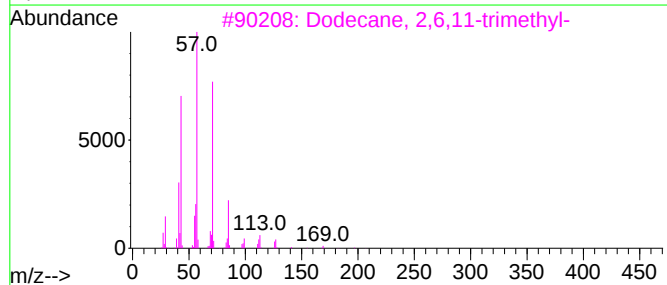
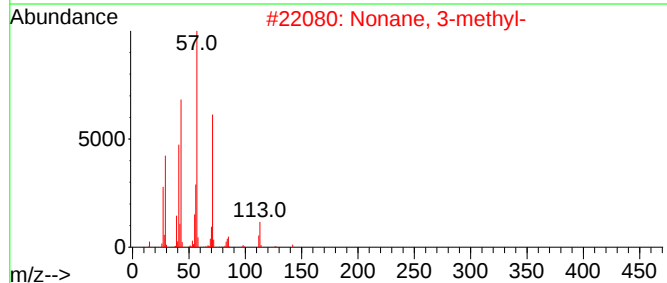
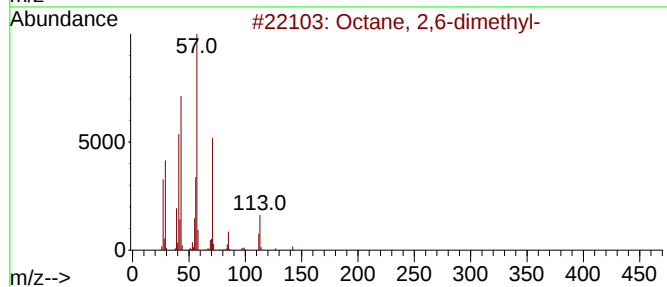
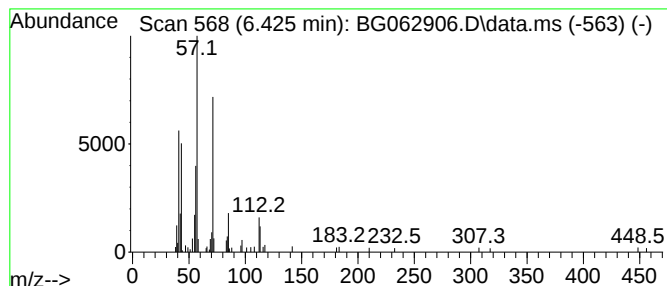
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.425	2.83 ng/ul	60772	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	80
2	Nonane, 3-methyl-			142	C10H22	005911-04-6	72
3	Dodecane, 2,6,11-trimethyl-			212	C15H32	031295-56-4	59
4	Butane, 2,2-dimethyl-			86	C6H14	000075-83-2	53
5	Octane, 3,5-dimethyl-			142	C10H22	015869-93-9	53



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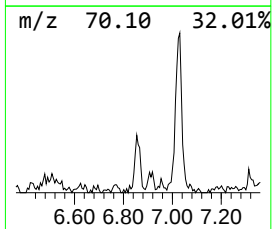
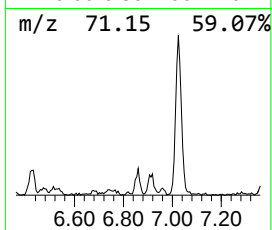
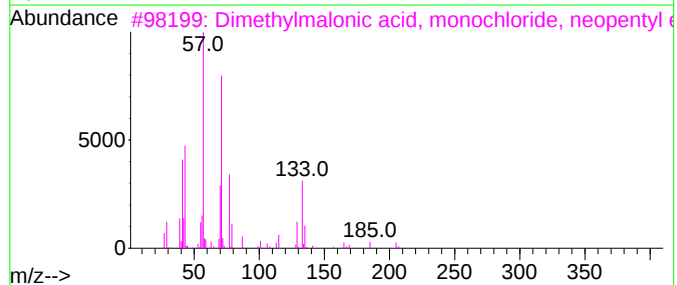
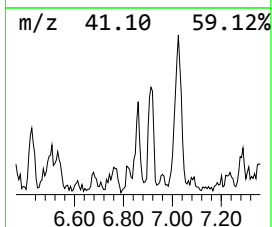
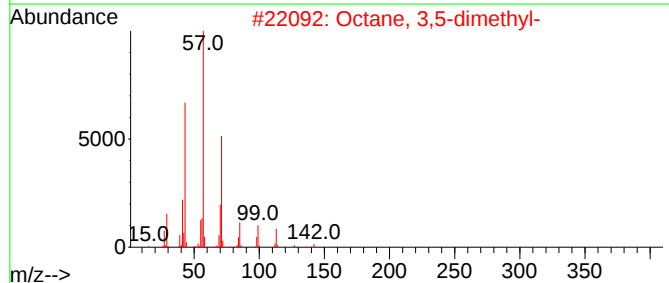
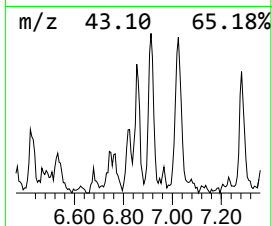
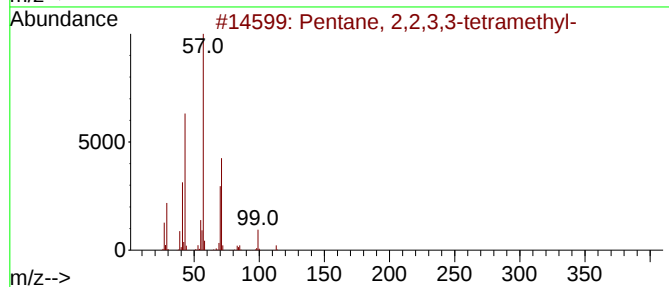
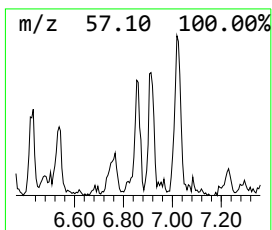
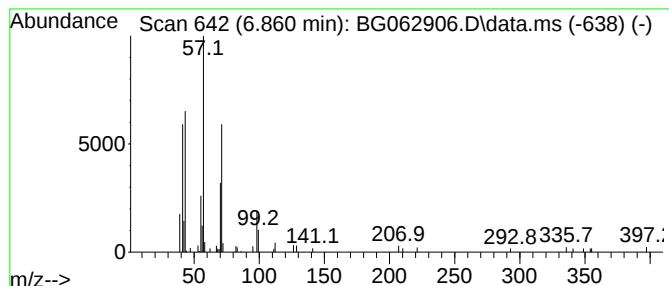
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.860	4.42 ng/ul	94879	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	64
2			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	59
3			Dimethylmalonic acid, monochlori...	220	C10H17ClO3	1000361-76-0	53
4			Tridecane, 6-methyl-	198	C14H30	013287-21-3	53
5			Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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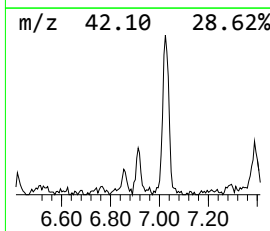
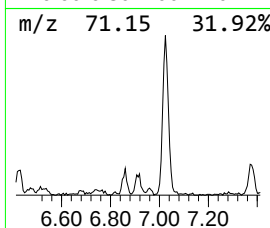
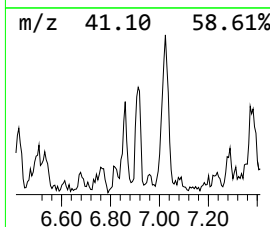
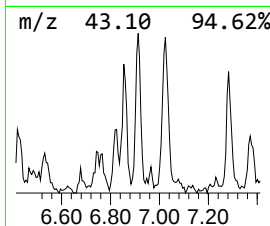
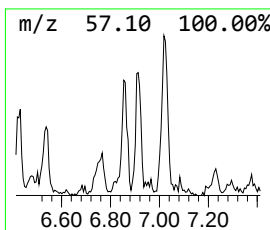
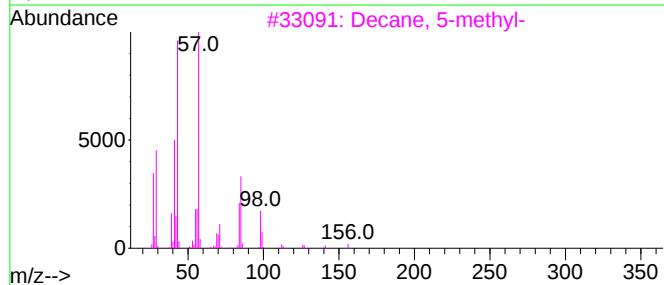
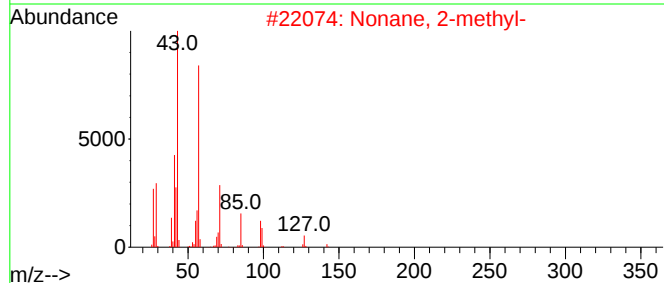
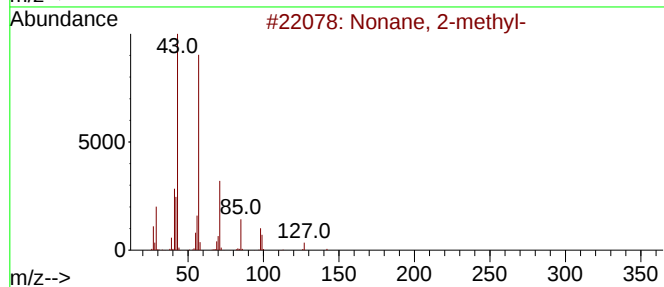
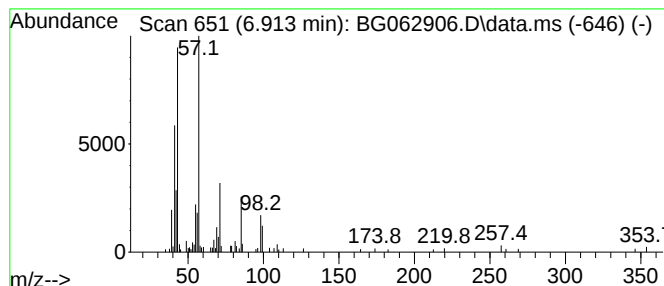
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.913	5.29 ng/ul	113592	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	64
2			Nonane, 2-methyl-	142	C10H22	000871-83-0	53
3			Decane, 5-methyl-	156	C11H24	013151-35-4	50
4			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	50
5			Nonane	128	C9H20	000111-84-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

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BNA_G
ClientSampleId :
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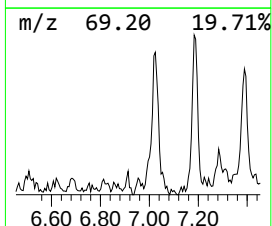
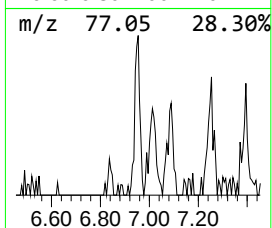
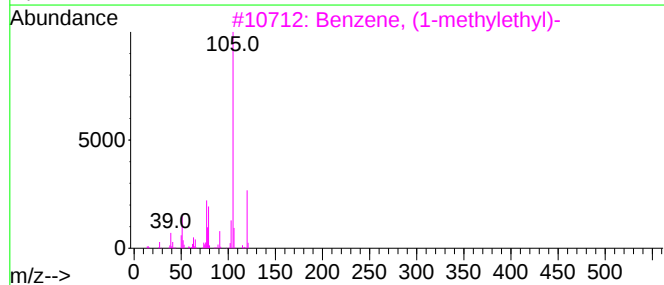
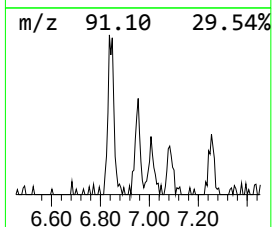
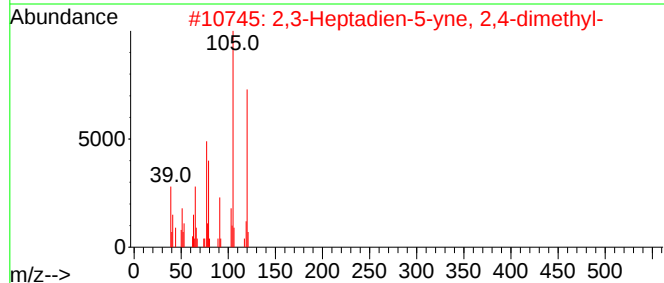
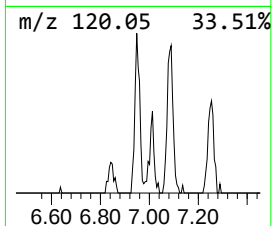
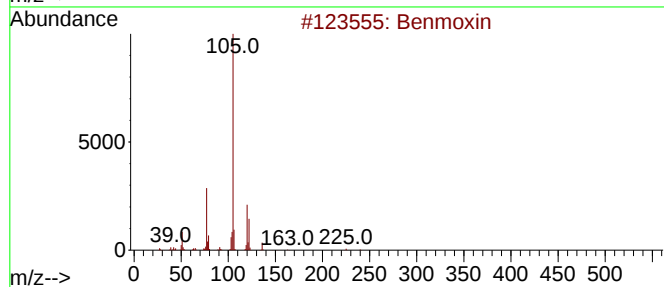
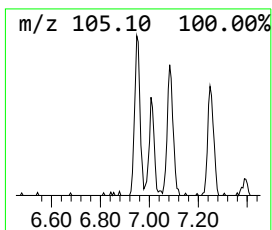
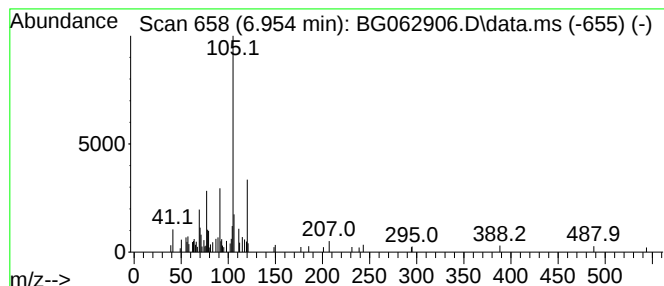
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benmoxin Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.954	3.30 ng/ul	70795	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benmoxin	240	C15H16N2O	007654-03-7	64
2			2,3-Heptadien-5-yne, 2,4-dimethyl-	120	C9H12	041898-89-9	59
3			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	58
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	58
5			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
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Operator : JU/RC
Sample : P3878-07ME
Misc :
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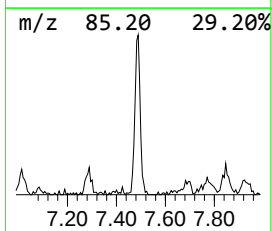
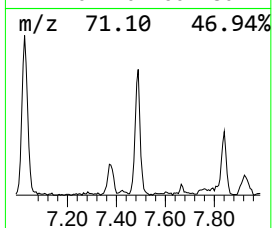
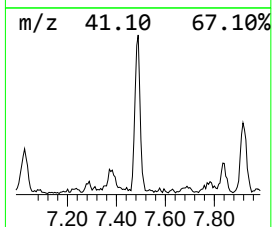
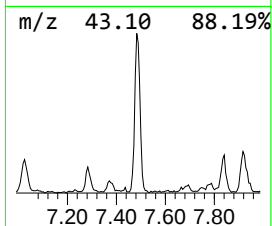
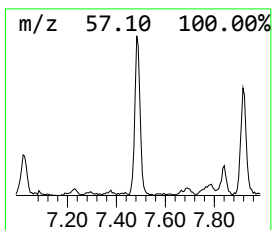
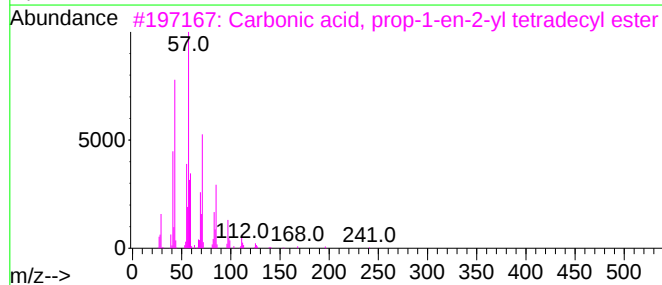
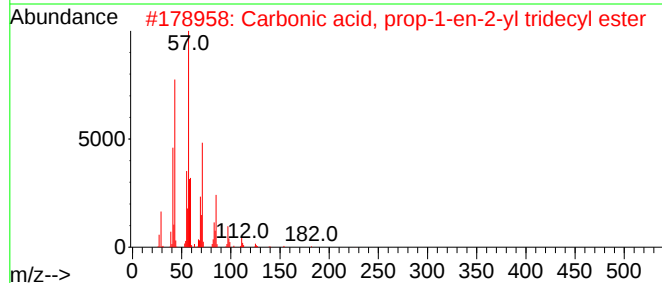
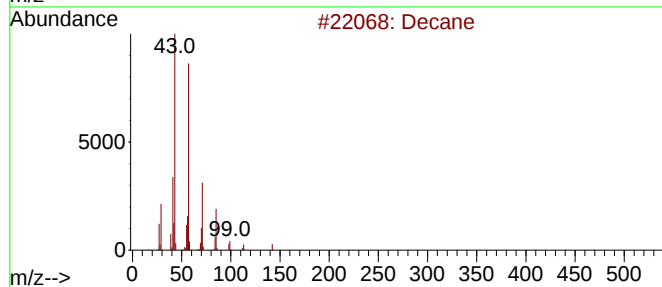
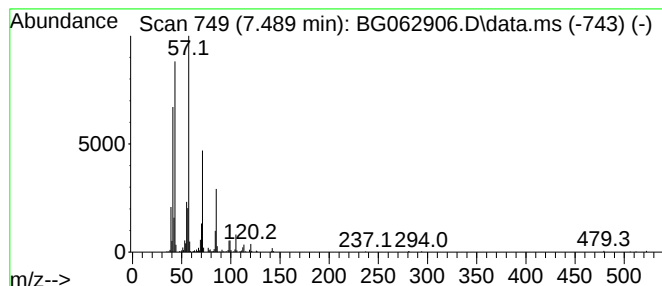
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.489	29.49 ng/ul	632780	1,4-Dichlorobenzene-d4	7.859	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	93
2	Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	72
3	Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	72
4	Undecane	156	C11H24	001120-21-4	64
5	Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
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Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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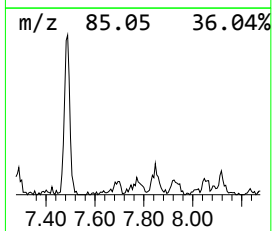
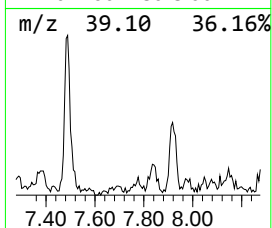
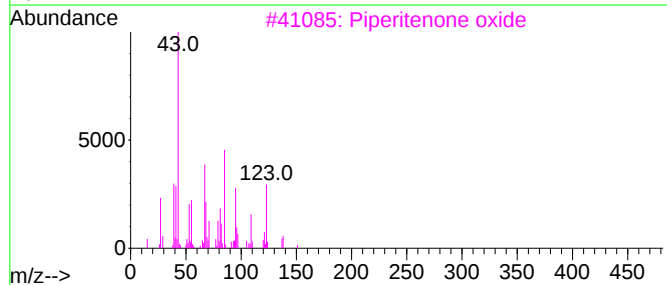
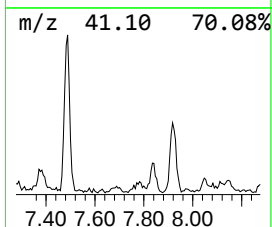
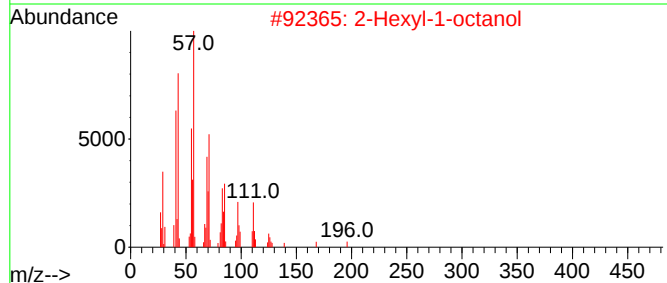
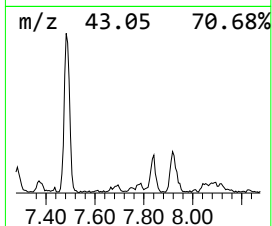
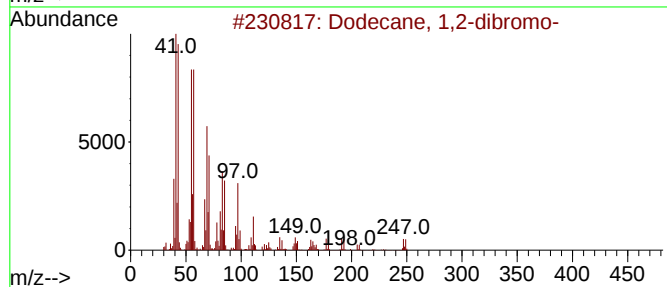
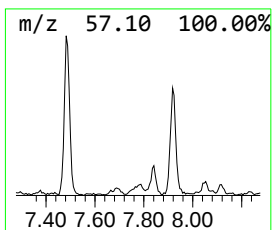
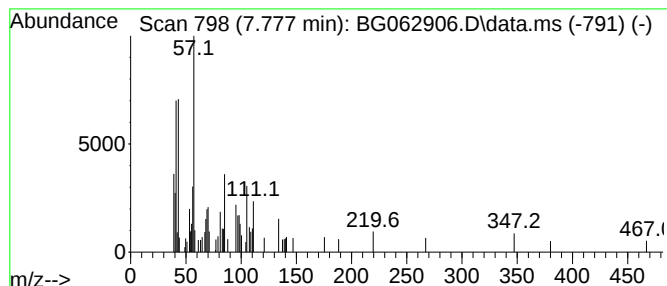
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.777	2.76 ng/ul	59254	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 1,2-dibromo-	326	C12H24Br2	055334-42-4	25
2			2-Hexyl-1-octanol	214	C14H30O	019780-79-1	22
3			Piperitenone oxide	166	C10H14O2	035178-55-3	22
4			3-(6,6-Dimethyl-5-oxohept-2-enyl...	250	C16H26O2	083040-97-5	22
5			1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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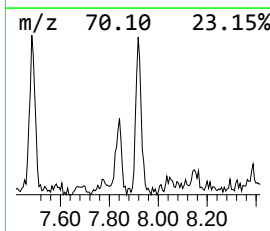
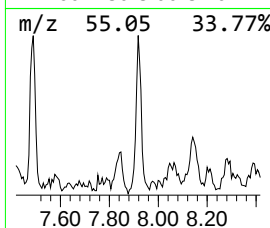
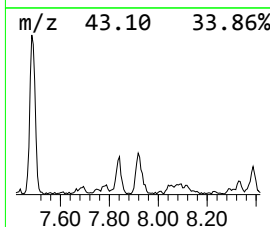
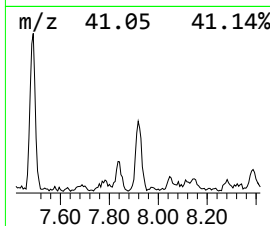
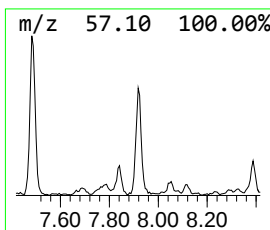
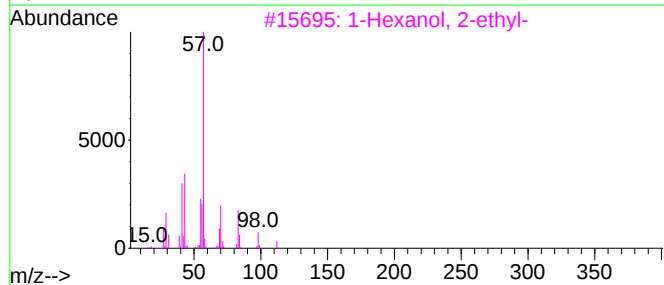
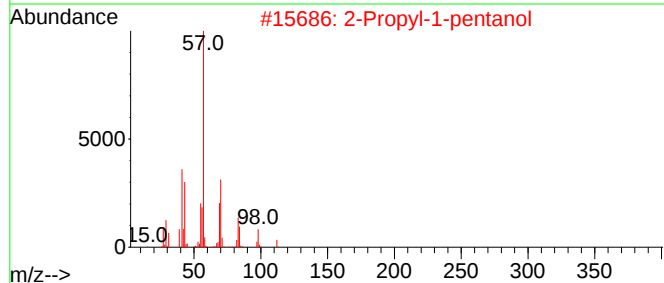
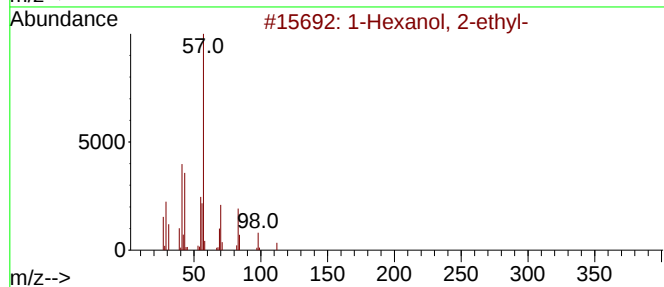
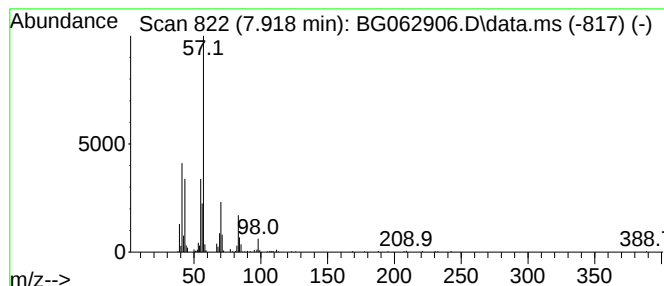
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Hexanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.918	11.73 ng/ul	251757	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
2	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	59
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	56
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	53
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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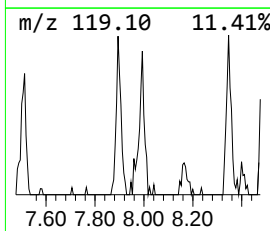
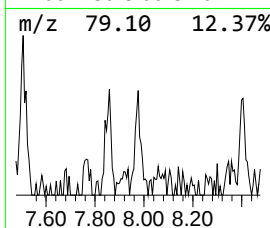
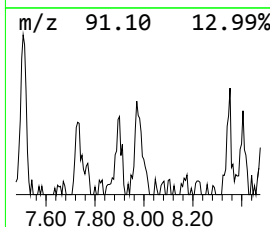
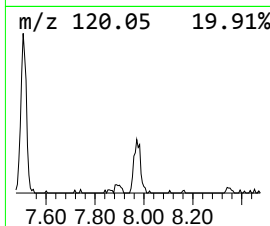
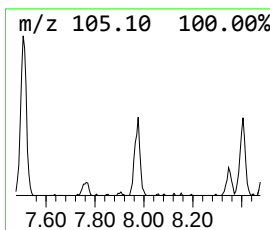
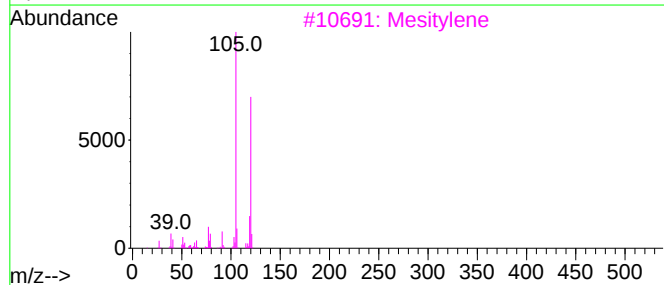
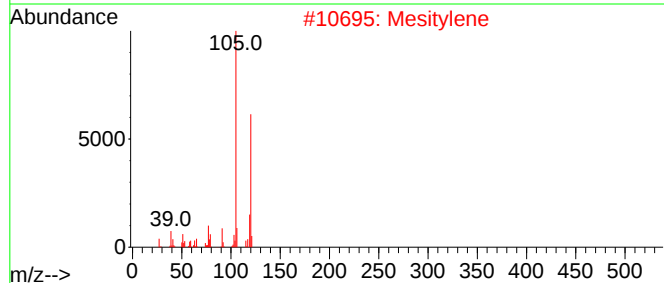
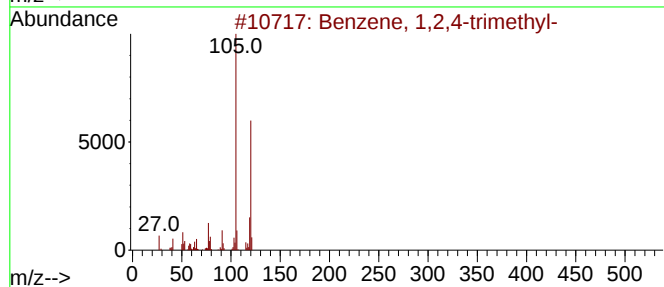
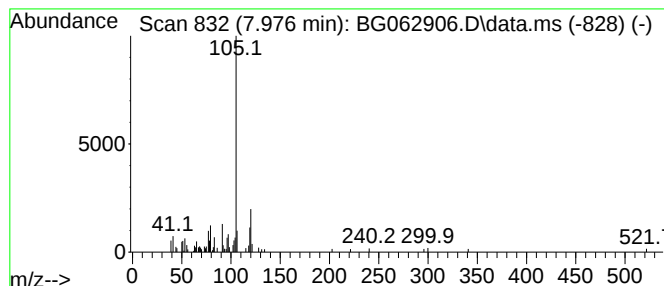
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,4-trimethyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.976	2.72 ng/ul	58290	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	72
2			Mesitylene	120	C9H12	000108-67-8	64
3			Mesitylene	120	C9H12	000108-67-8	64
4			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
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Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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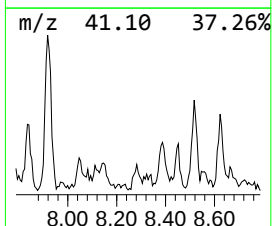
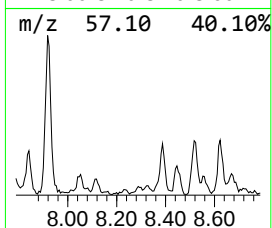
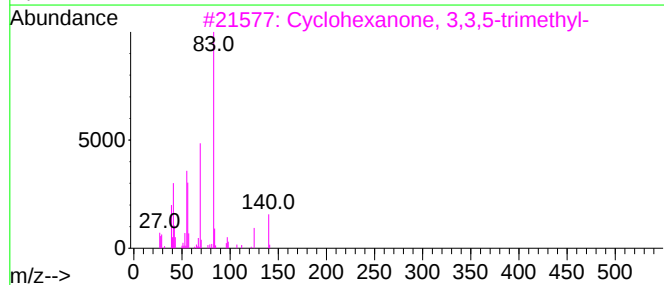
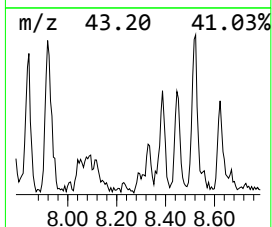
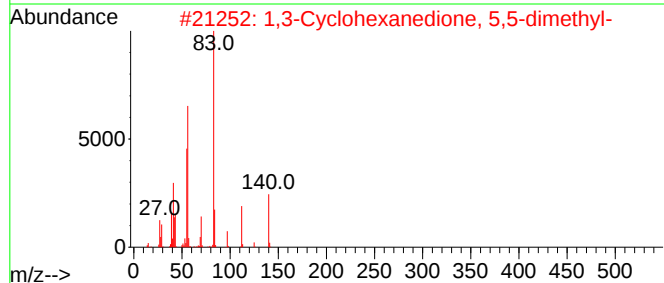
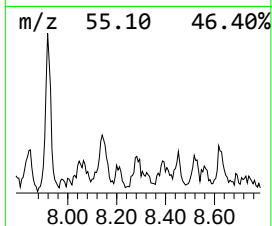
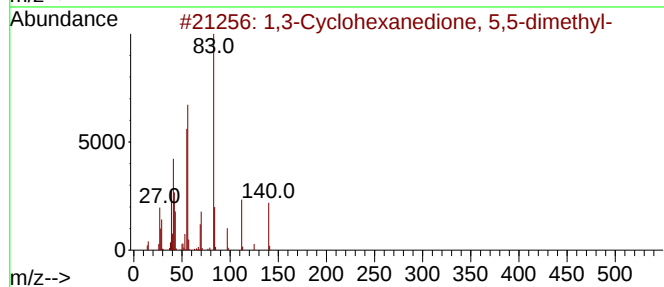
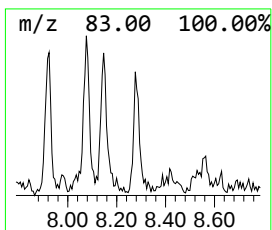
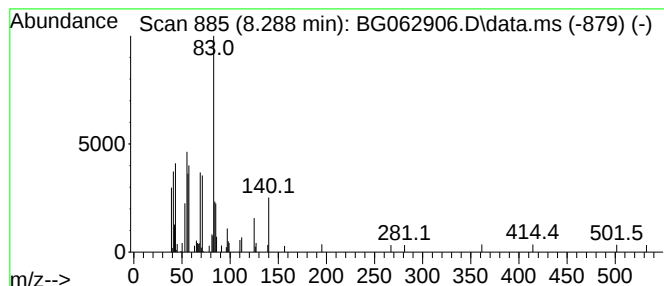
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,3-Cyclohexanedione, 5,5-d... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.288	2.99 ng/ul	64128	1,4-Dichlorobenzene-d4			7.859

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	59
2		1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	59
3		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	53
4		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	53
5		Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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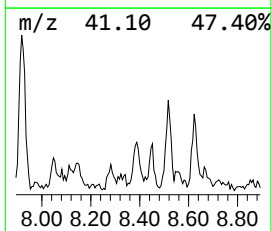
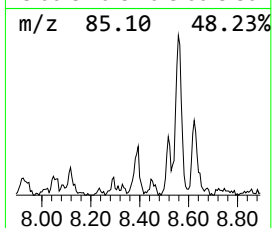
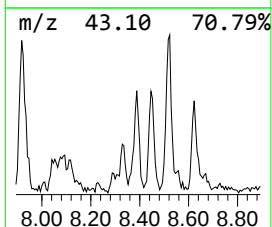
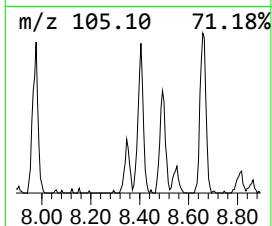
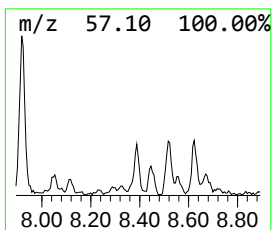
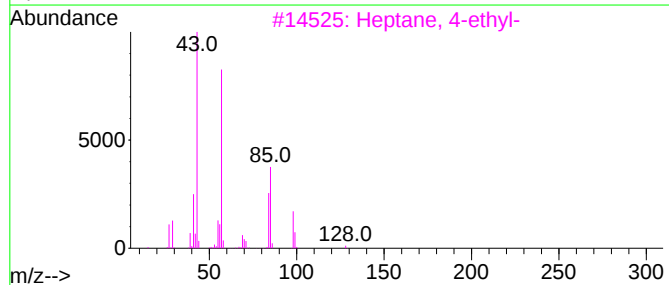
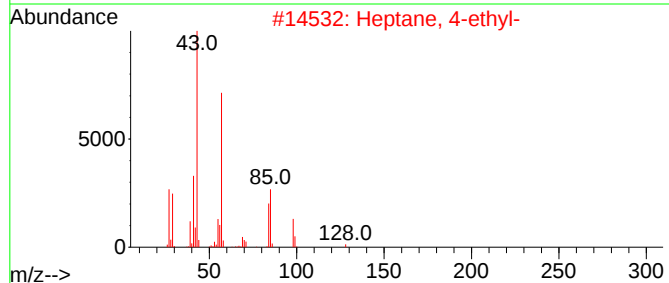
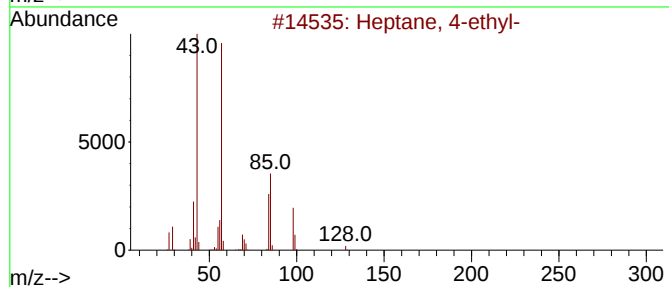
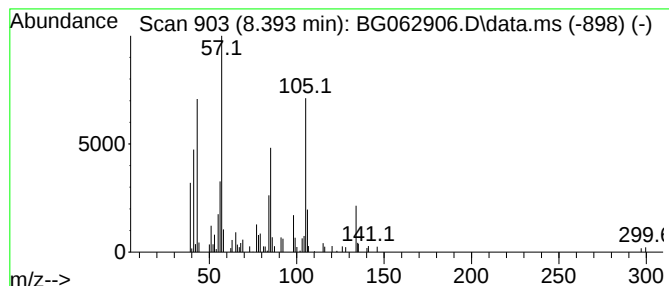
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.393	5.71 ng/ul	122533	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 4-ethyl-	128	C9H20	002216-32-2	47
2			Heptane, 4-ethyl-	128	C9H20	002216-32-2	47
3			Heptane, 4-ethyl-	128	C9H20	002216-32-2	43
4			Oxalic acid, heptyl isohexyl ester	272	C15H28O4	1000309-33-1	43
5			Hydratropic acid, isohexyl ester	234	C15H22O2	1000415-06-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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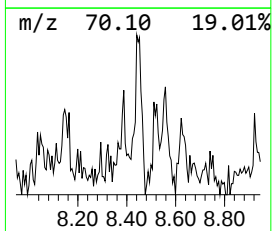
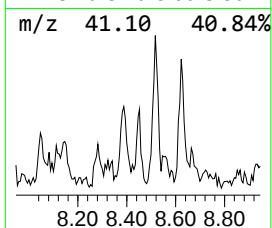
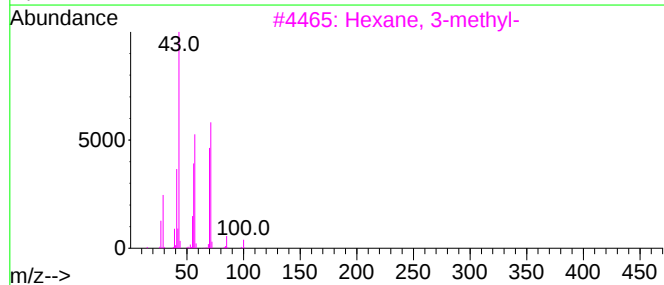
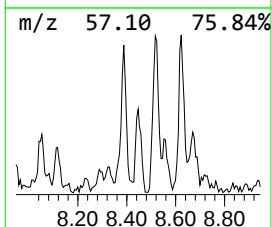
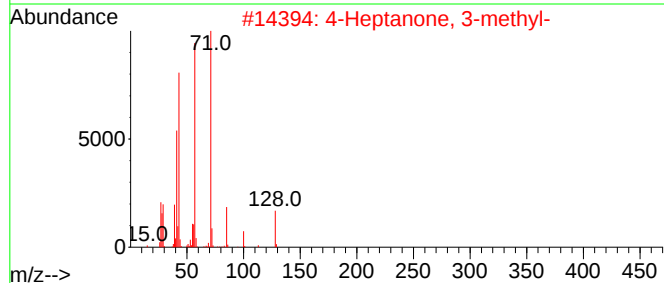
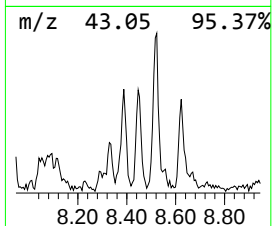
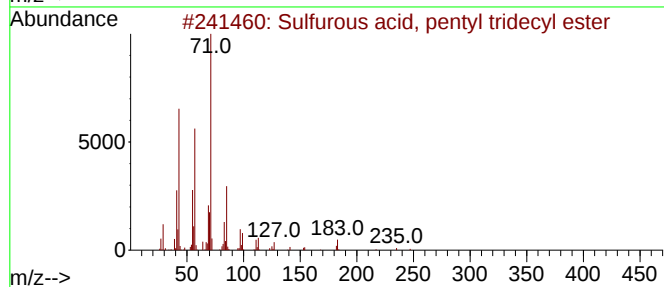
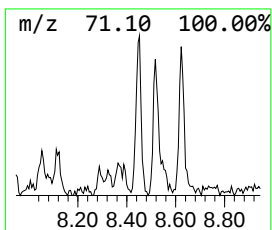
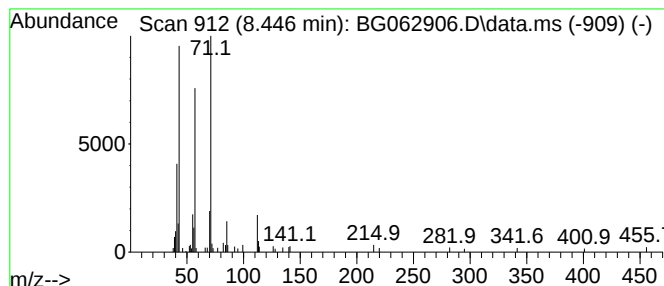
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Sulfurous acid, pentyl trid... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.446	2.65 ng/ul	56836	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, pentyl tridecyl ...	334	C18H38O3S	1000309-14-6	64
2			4-Heptanone, 3-methyl-	128	C8H16O	015726-15-5	64
3			Hexane, 3-methyl-	100	C7H16	000589-34-4	64
4			3-Octen-2-ol	128	C8H16O	076649-14-4	59
5			Nonane, 1-iodo-	254	C9H19I	004282-42-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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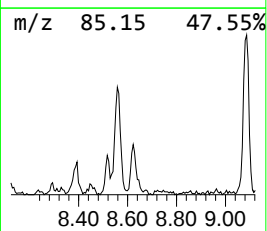
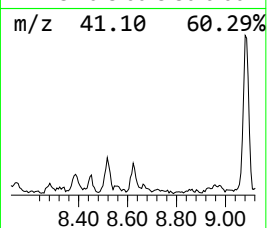
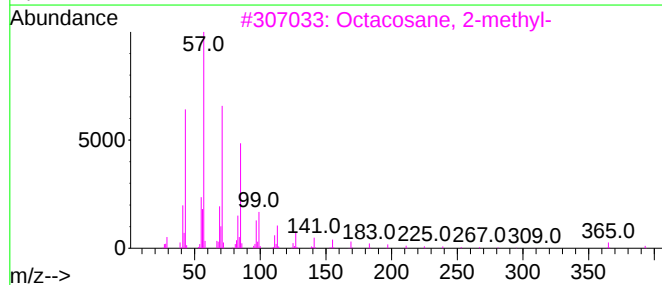
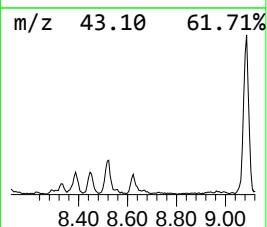
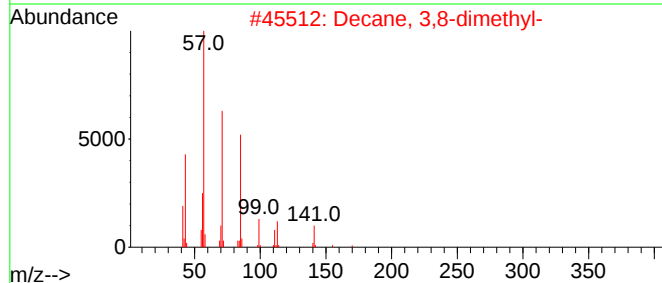
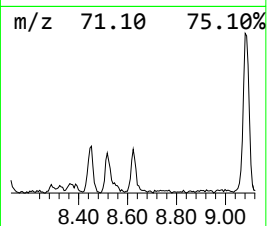
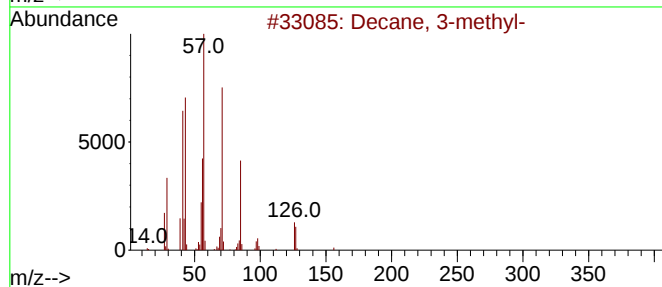
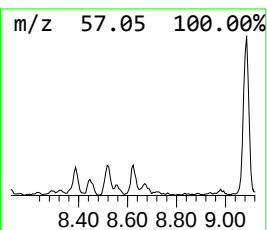
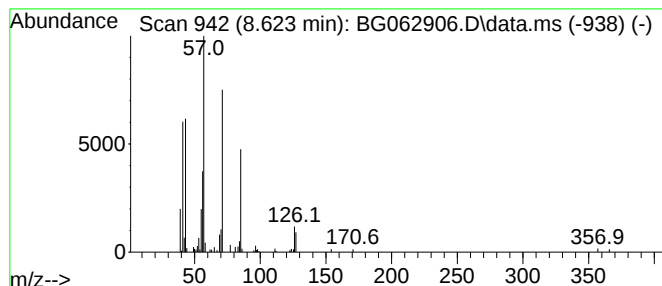
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.623	4.35 ng/ul	93367	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	83
2			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	72
4			Dodecane, 3-methyl-	184	C13H28	017312-57-1	72
5			Octane, 2-methyl-	128	C9H20	003221-61-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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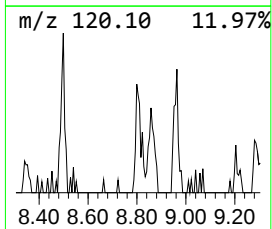
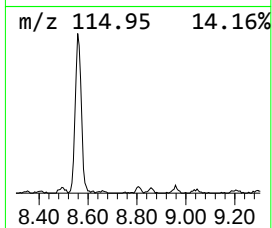
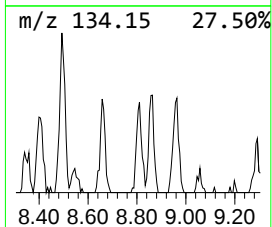
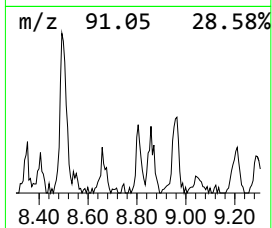
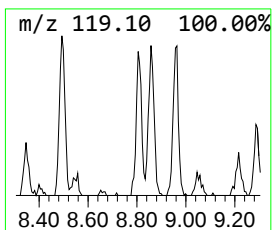
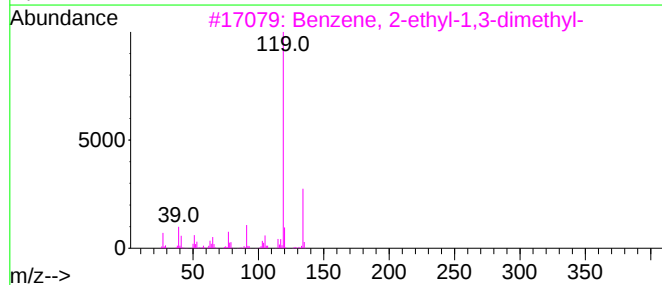
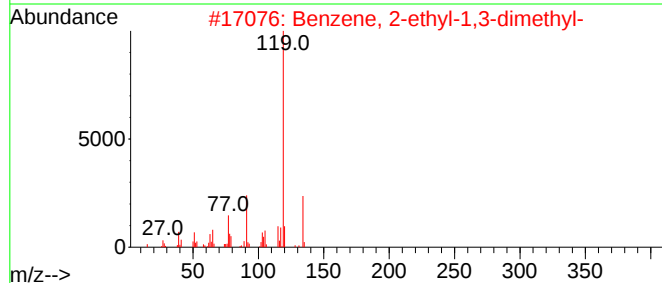
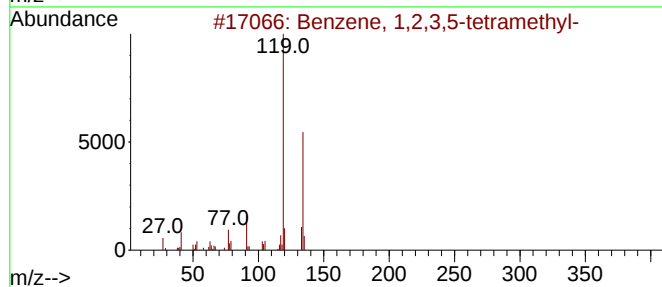
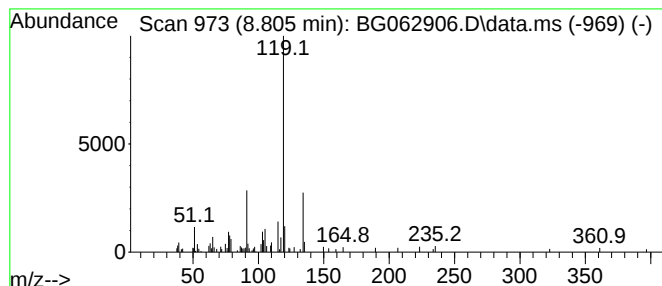
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.805	2.56 ng/ul	54998	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87
2			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	87
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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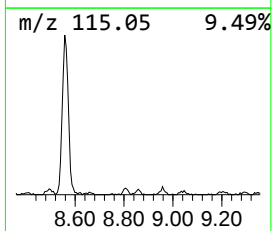
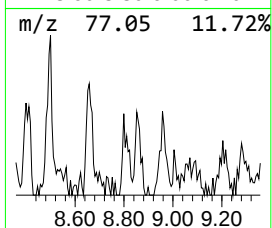
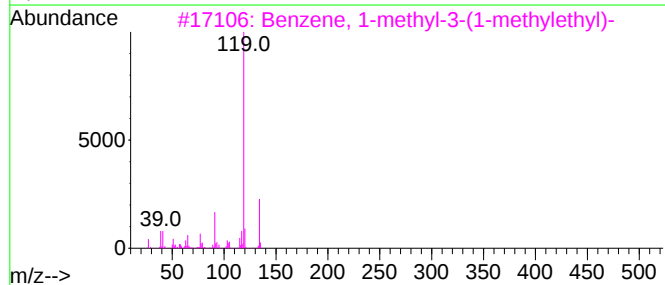
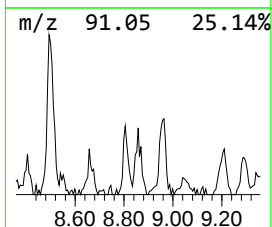
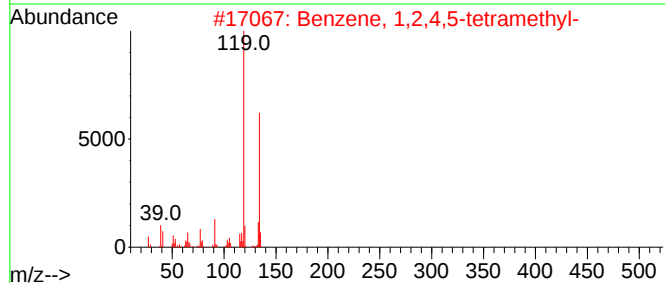
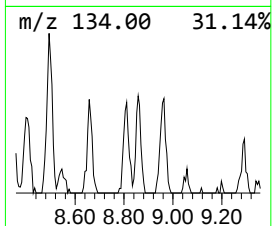
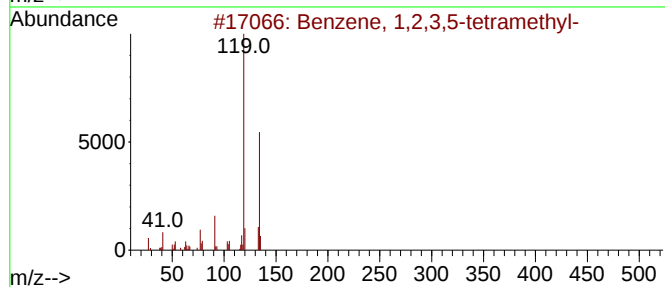
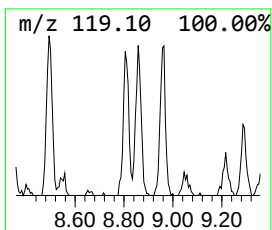
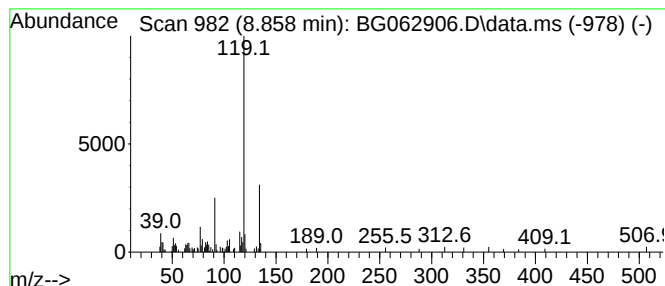
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.858	3.37 ng/ul	72354	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81
3			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	81
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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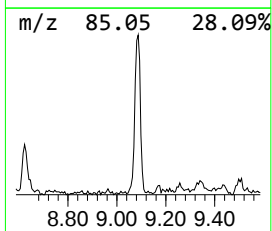
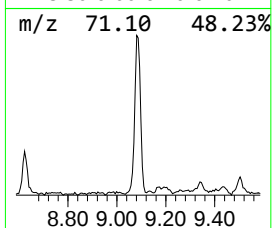
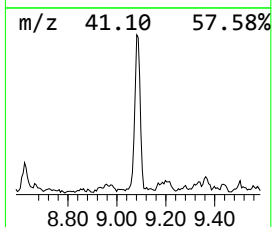
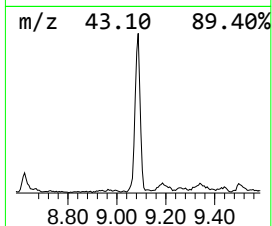
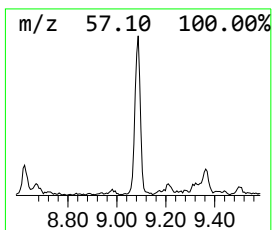
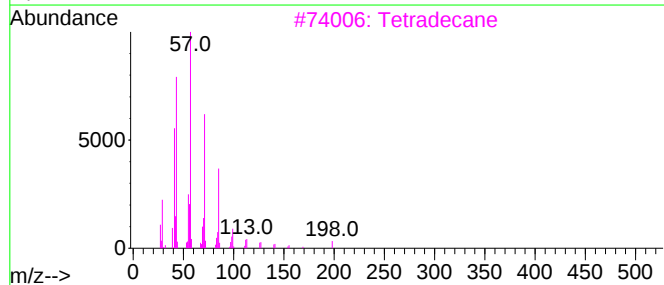
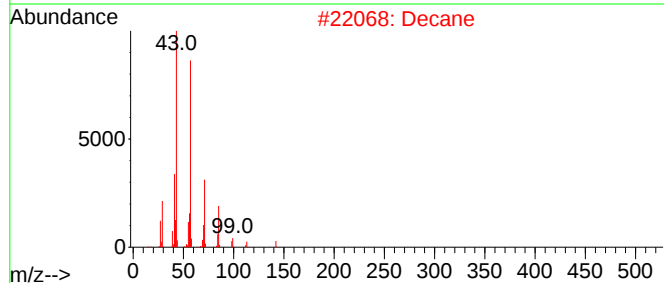
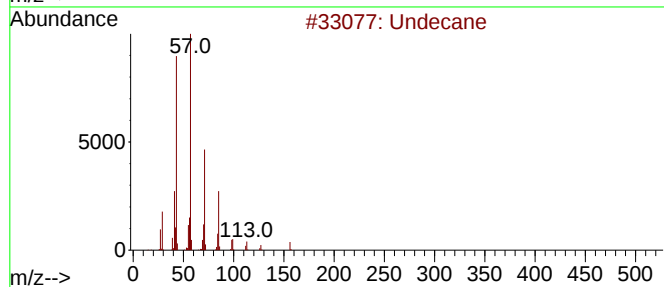
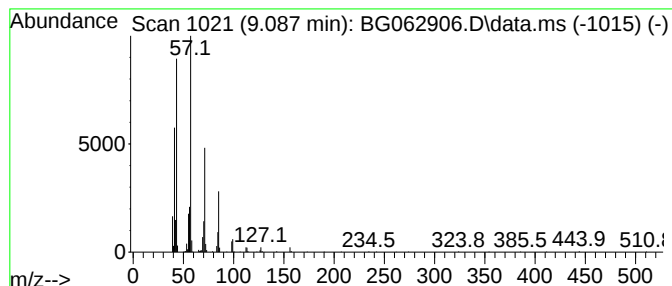
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.087	29.10 ng/ul	624369	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Decane	142	C10H22	000124-18-5	90
3			Tetradecane	198	C14H30	000629-59-4	90
4			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	86
5			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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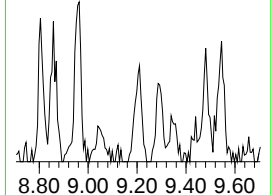
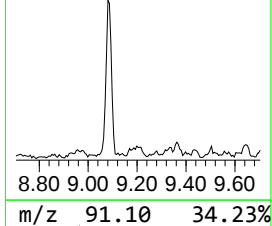
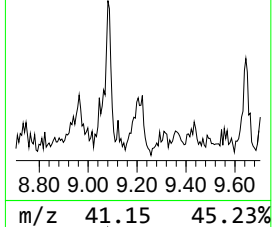
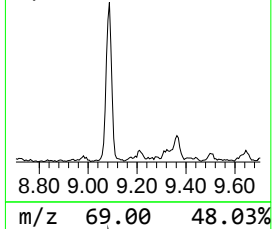
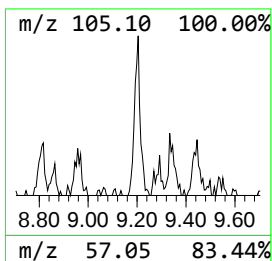
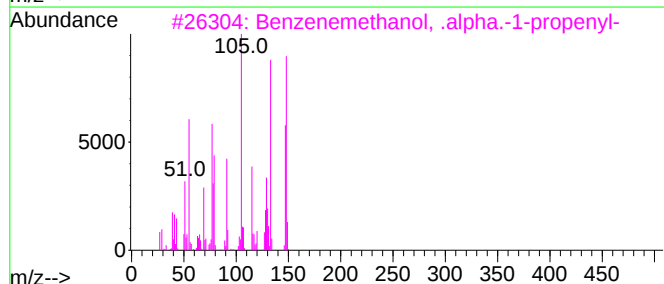
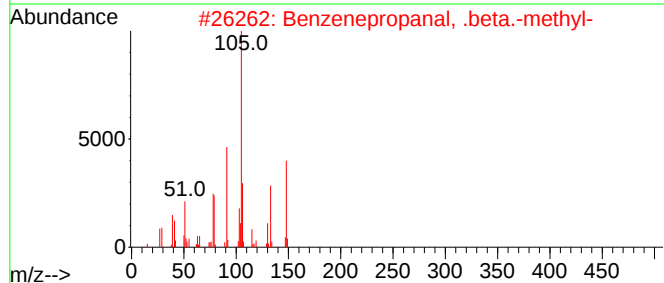
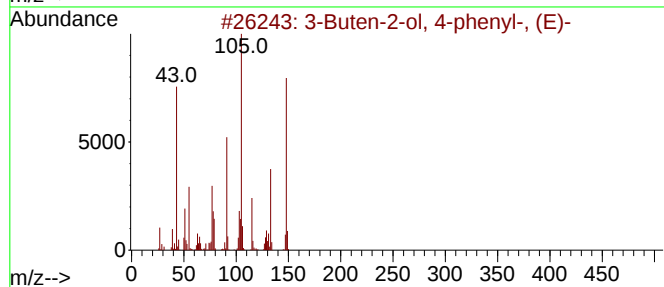
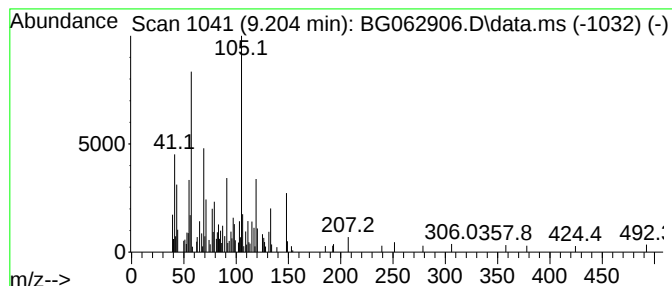
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.204	7.00 ng/ul	150250	1,4-Dichlorobenzene-d4	7.859

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	38
2			Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	30
3			Benzenemethanol, .alpha.-1-prope...	148	C10H12O	003347-57-7	30
4			Tricyclo[4.2.1.1(2,5)]deca-3,7-d...	176	C11H12O2	070220-88-1	27
5			3-Buten-2-ol, 4-phenyl-	148	C10H12O	017488-65-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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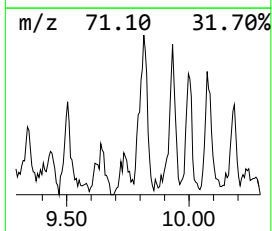
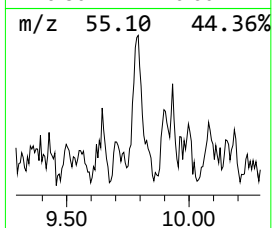
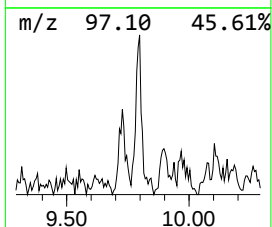
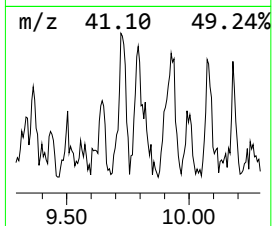
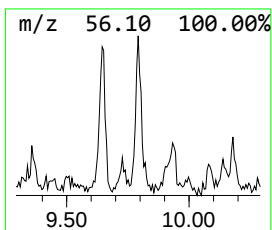
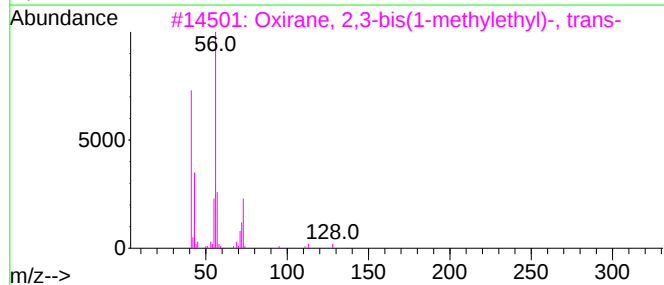
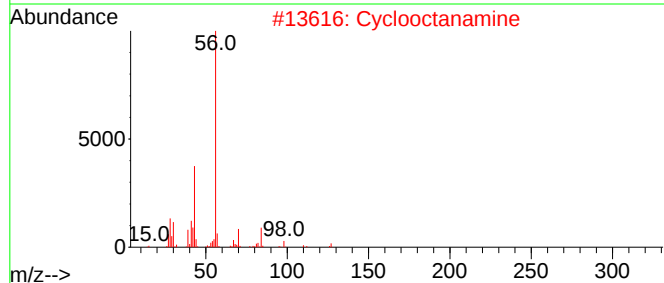
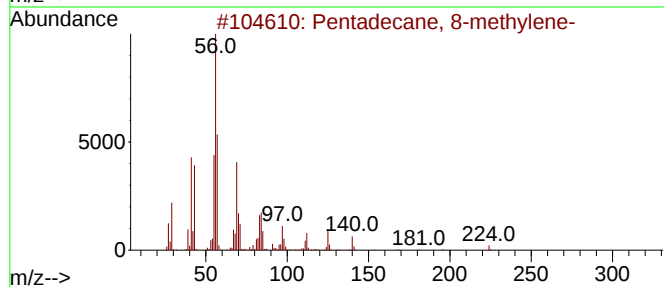
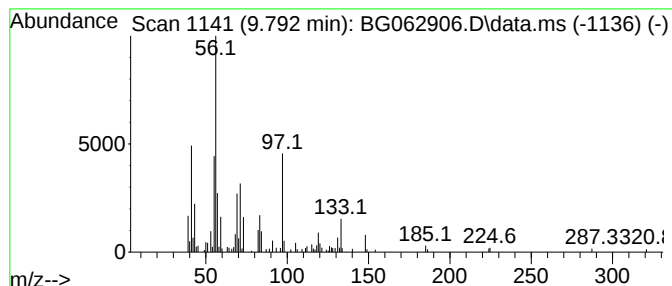
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.792	2.48 ng/ul	129929	Naphthalene-d8	10.644

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 8-methylene-	224	C16H32	055668-09-2	38
2		Cyclooctanamine	127	C8H17N	005452-37-9	35
3		Oxirane, 2,3-bis(1-methylethyl)-...	128	C8H16O	054644-32-5	35
4		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	35
5		1-Decene, 9-methyl-	154	C11H22	061142-78-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

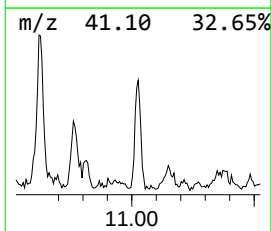
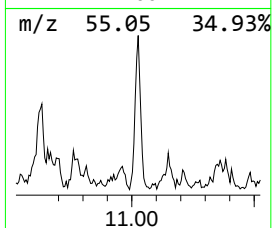
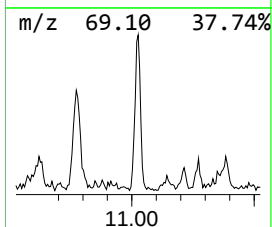
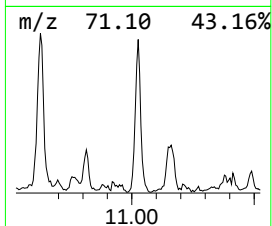
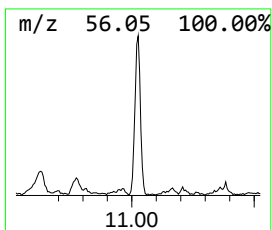
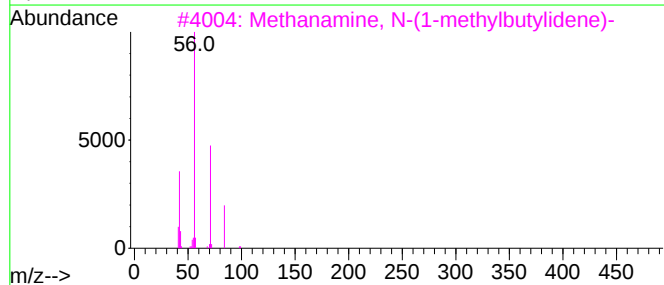
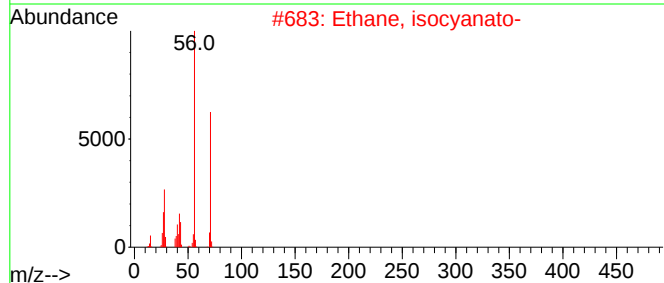
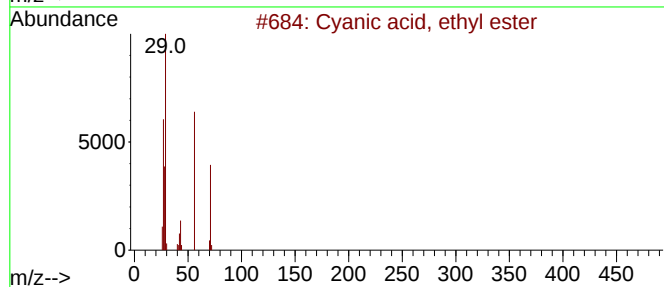
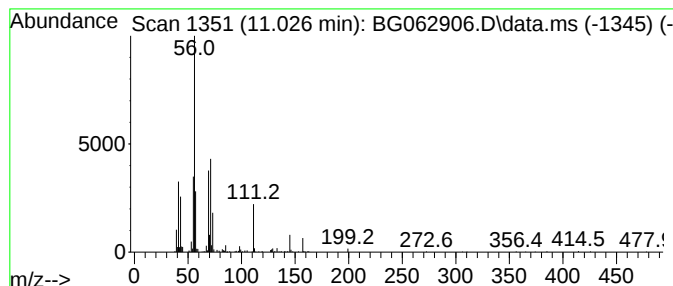
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 19 Cyanic acid, ethyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.026	8.93 ng/ul	468724	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyanic acid, ethyl ester	71	C3H5NO	000627-48-5	58
2	Ethane, isocyanato-	71	C3H5NO	000109-90-0	42
3	Methanamine, N-(1-methylbutylidene...	99	C6H13N	022431-09-0	42
4	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	38
5	Cyclohexylamine	99	C6H13N	000108-91-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

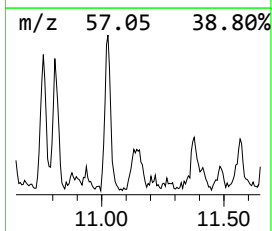
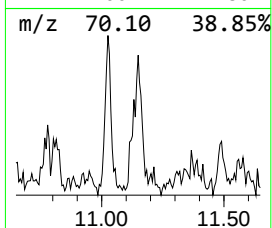
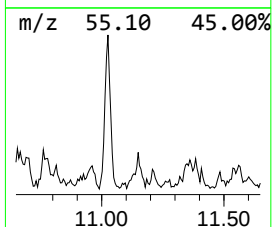
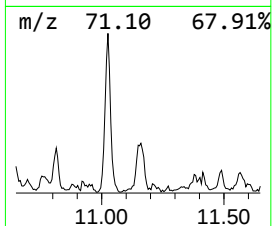
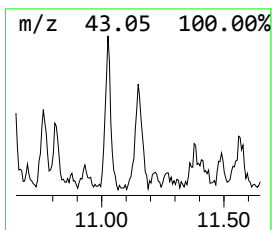
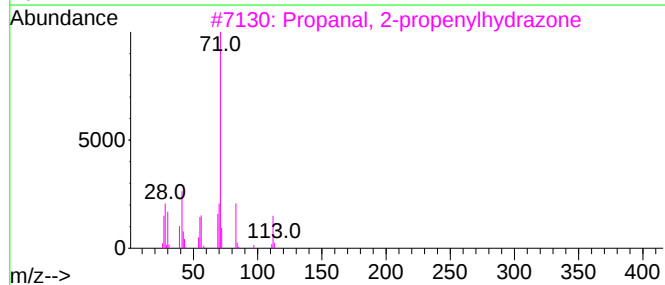
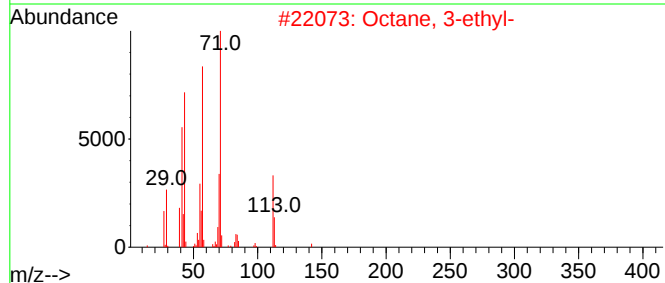
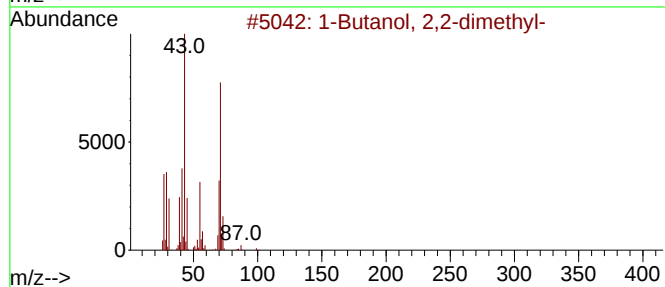
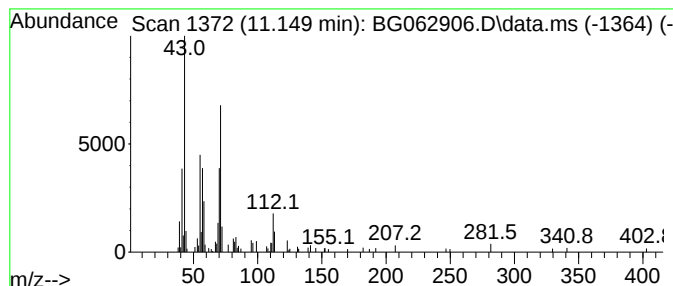
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 20 unknown-03 Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.149	2.87 ng/ul	150568	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Butanol, 2,2-dimethyl-	102	C6H14O	001185-33-7	47
2	Octane, 3-ethyl-	142	C10H22	005881-17-4	47
3	Propanal, 2-propenylhydrazone	112	C6H12N2	019031-78-8	43
4	Oxalic acid, 2-ethylhexyl pentyl...	272	C15H28O4	1000309-38-7	42
5	Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
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Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

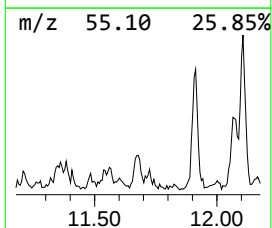
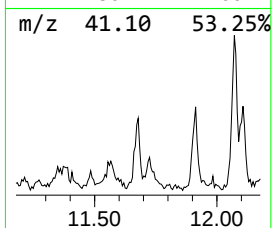
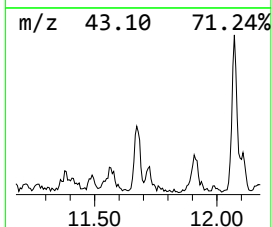
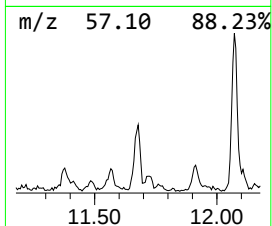
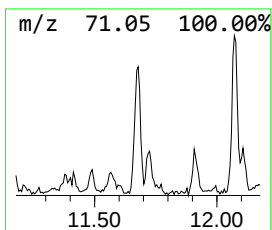
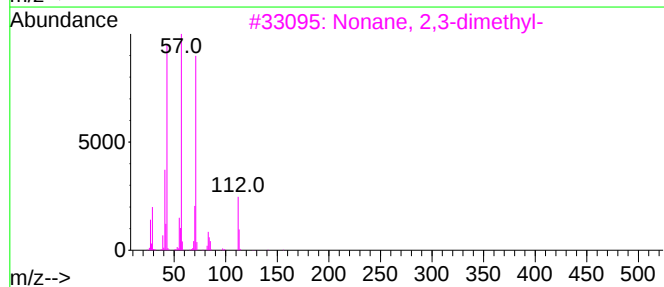
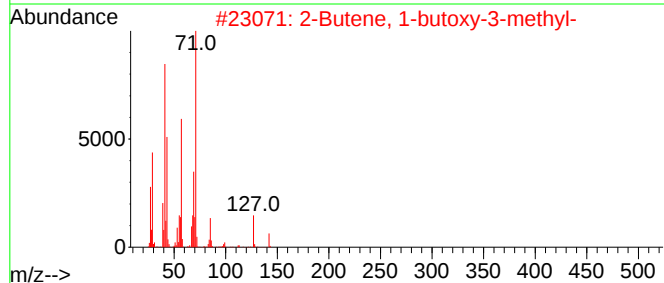
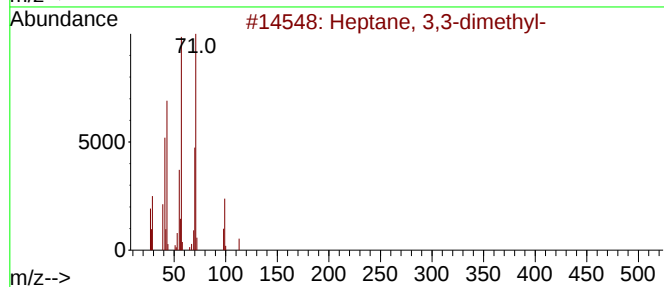
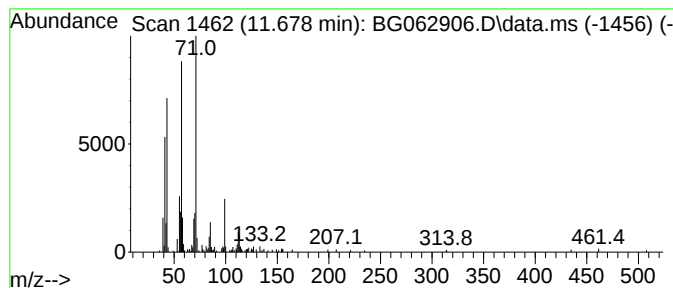
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.678	3.99 ng/ul	209570	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	59
2	2-Butene, 1-butoxy-3-methyl-	142	C9H18O	022094-02-6	53
3	Nonane, 2,3-dimethyl-	156	C11H24	002884-06-2	53
4	Octane, 2-bromo-	192	C8H17Br	000557-35-7	50
5	3-Bromooctane	192	C8H17Br	000999-64-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

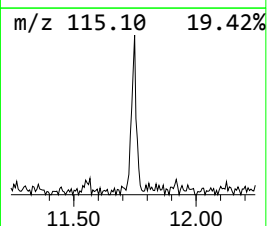
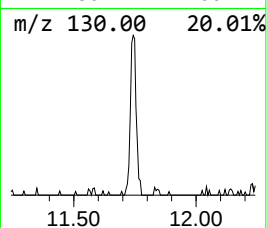
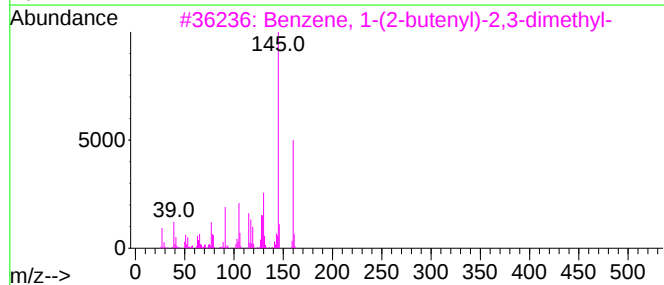
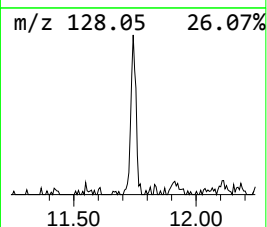
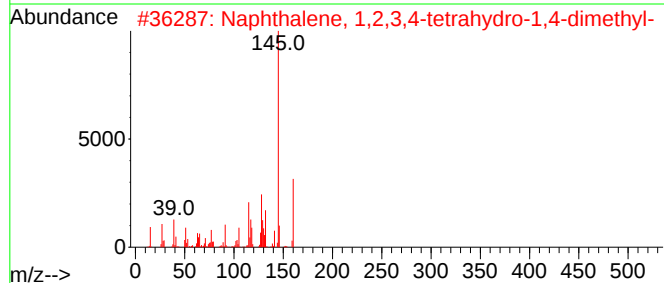
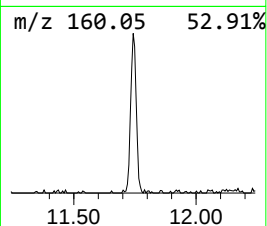
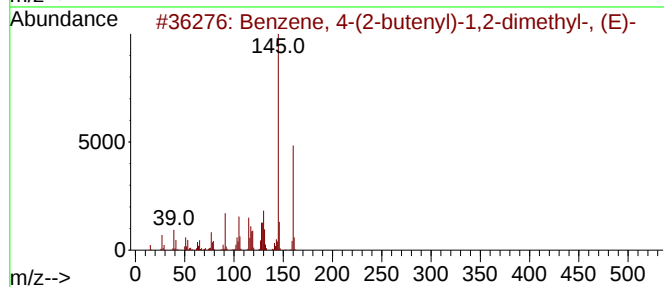
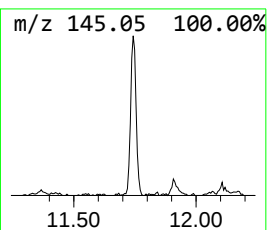
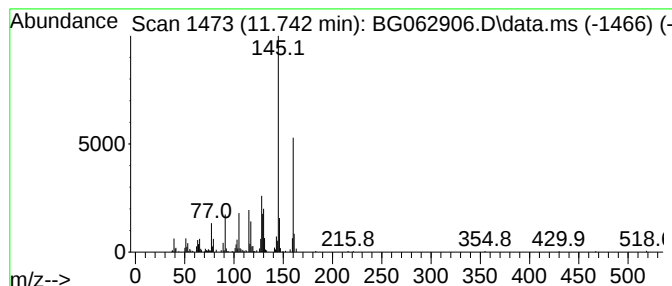
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.743	7.33 ng/ul	384398	Naphthalene-d8	10.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	90
2			Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	90
3			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	90
4			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	90
5			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
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Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

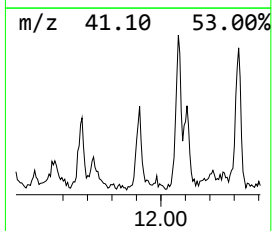
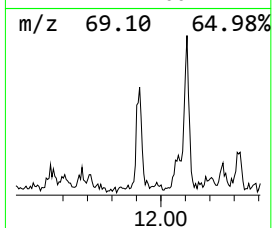
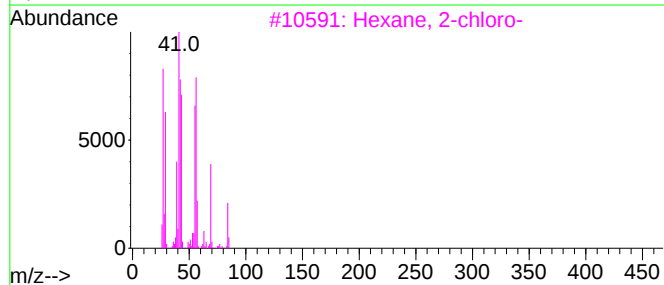
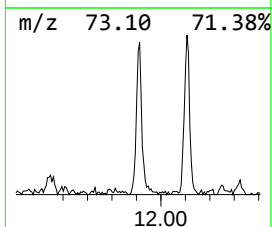
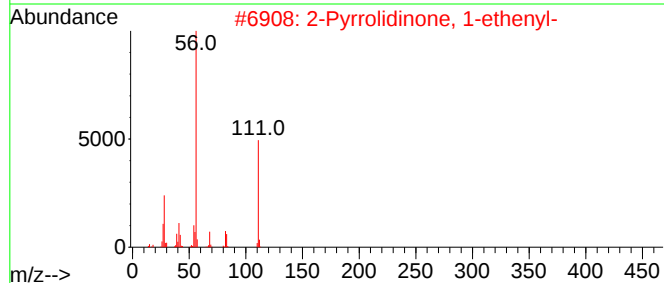
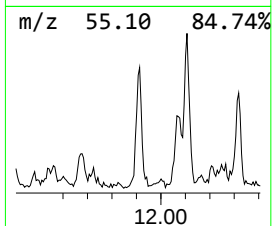
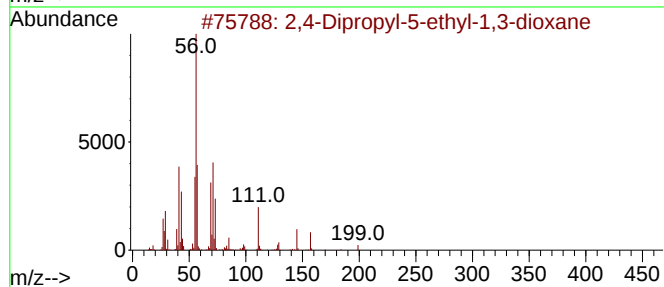
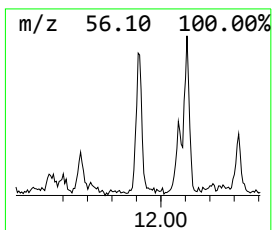
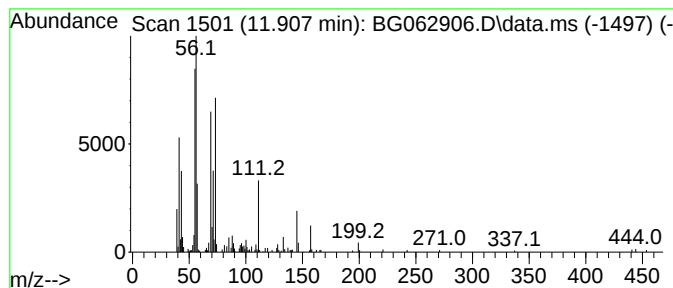
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 23 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.907	4.62 ng/ul	242336	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	59
2	2-Pyrrolidinone, 1-ethenyl-	111	C6H9NO	000088-12-0	43
3	Hexane, 2-chloro-	120	C6H13Cl	000638-28-8	38
4	1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	38
5	Ethanol, 2-(dodecyloxy)-	230	C14H30O2	004536-30-5	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

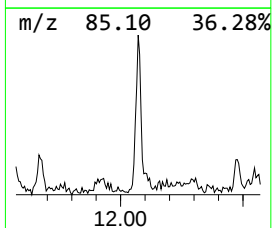
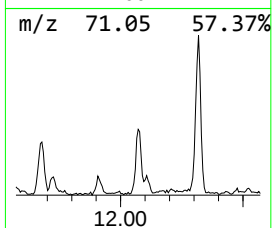
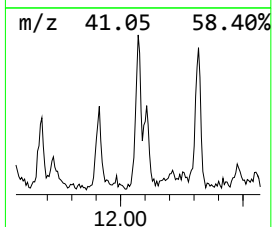
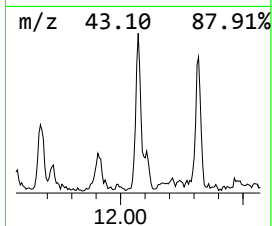
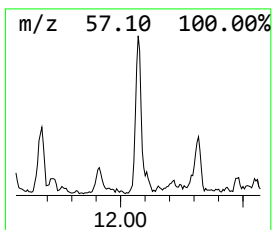
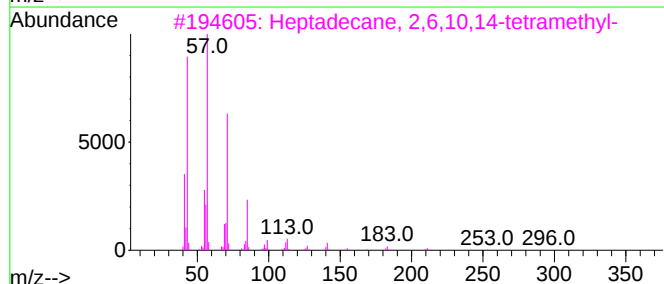
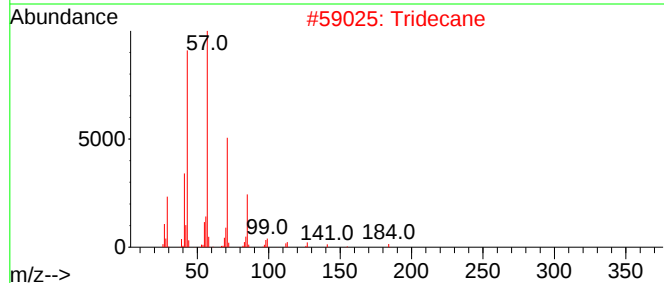
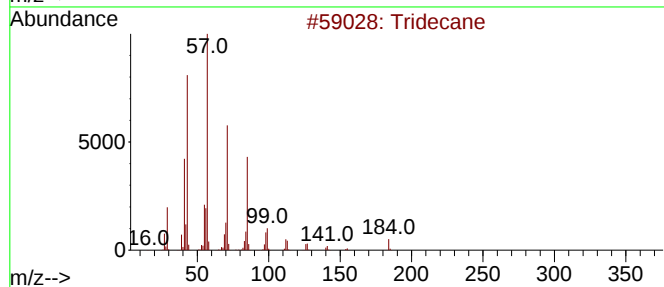
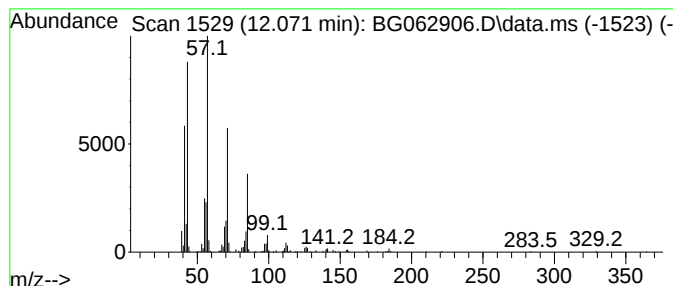
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.072	7.14 ng/ul	374688	Naphthalene-d8	10.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Tridecane	184	C13H28	000629-50-5	87
3			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	86
4			Hexadecane	226	C16H34	000544-76-3	83
5			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

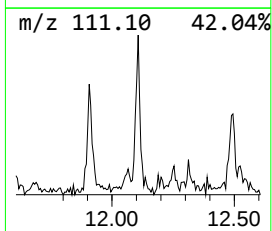
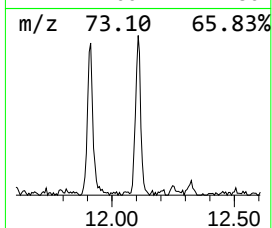
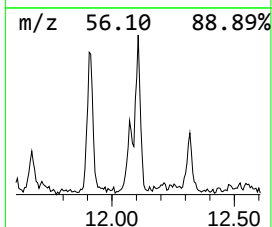
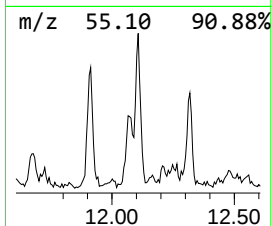
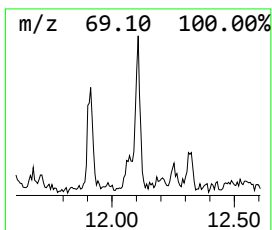
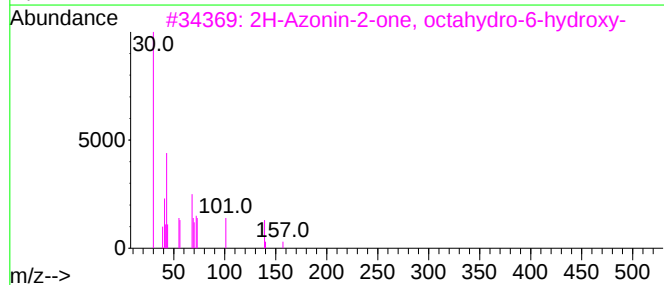
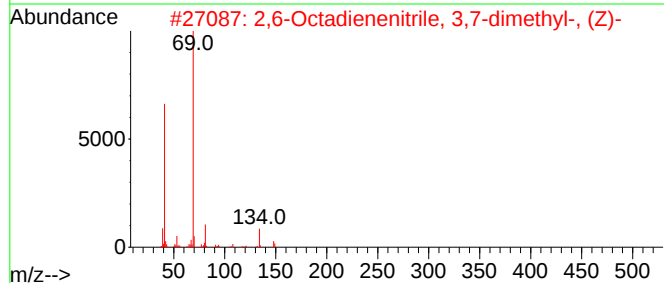
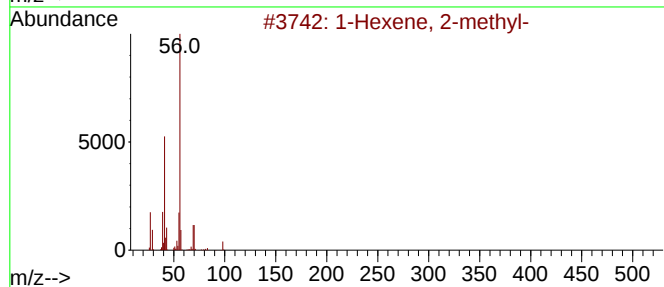
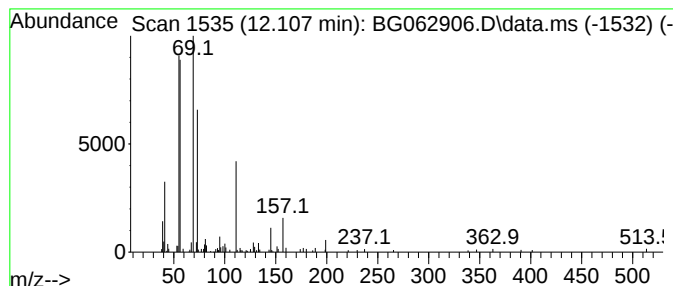
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.107	4.61 ng/ul	241658	Naphthalene-d8	10.644

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	10	
2	2,6-Octadienenitrile, 3,7-dimeth...	149	C10H15N	031983-27-4	10	
3	2H-Azonin-2-one, octahydro-6-hyd...	157	C8H15NO2	023435-07-6	9	
4	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	9	
5	Succinic anhydride	100	C4H4O3	000108-30-5	9	



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

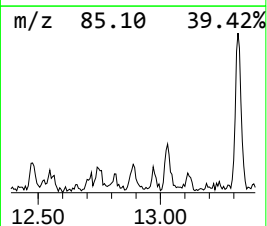
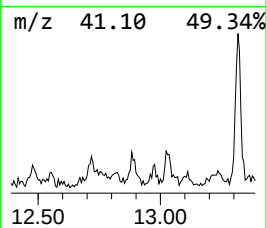
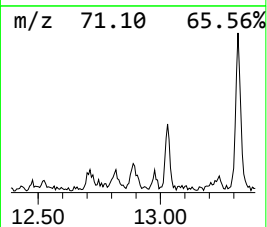
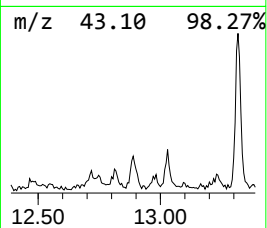
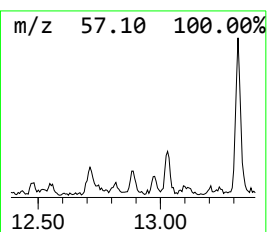
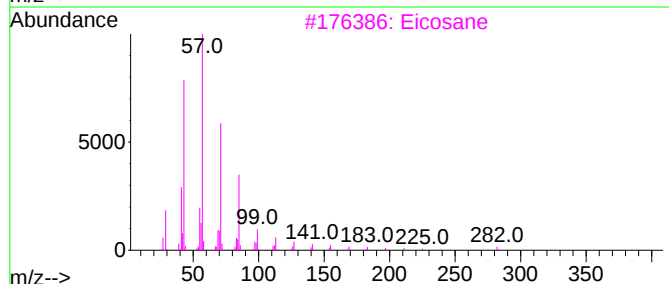
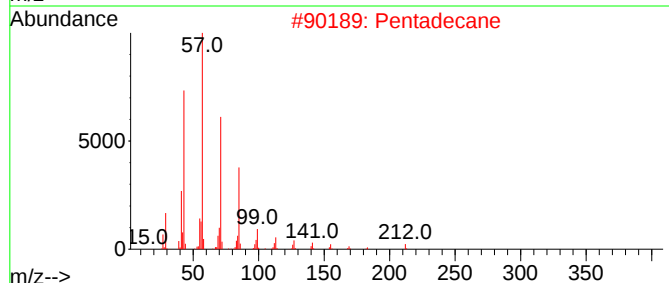
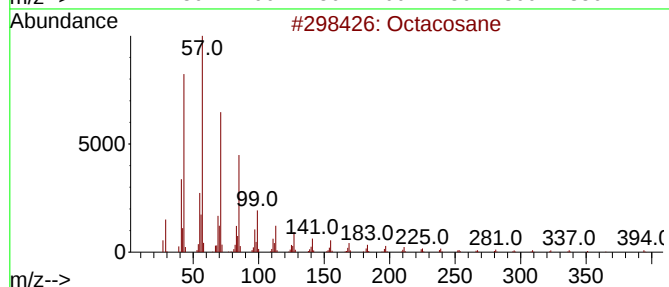
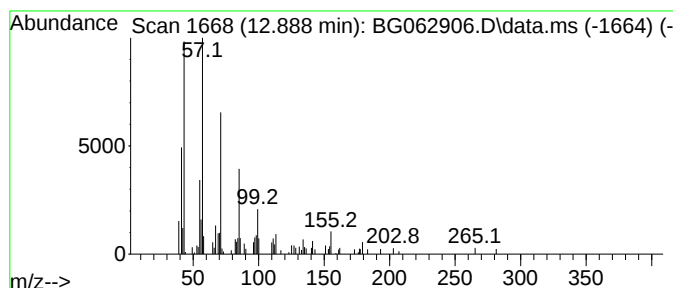
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.888	2.29 ng/ul	91168	Acenaphthene-d10		14.486	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	80
2	Pentadecane		212	C15H32	000629-62-9	72
3	Eicosane		282	C20H42	000112-95-8	72
4	Pentadecane		212	C15H32	000629-62-9	72
5	Heptadecane		240	C17H36	000629-78-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

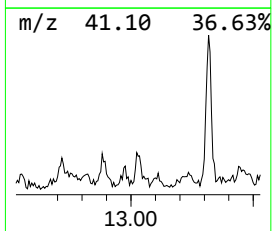
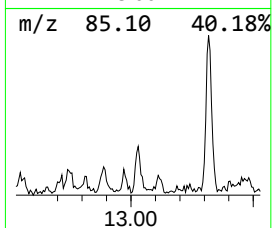
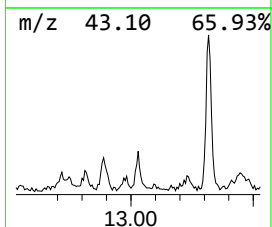
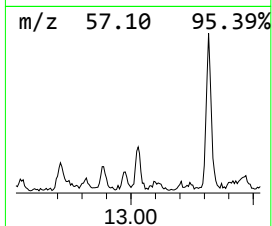
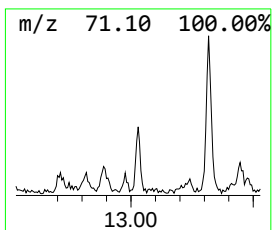
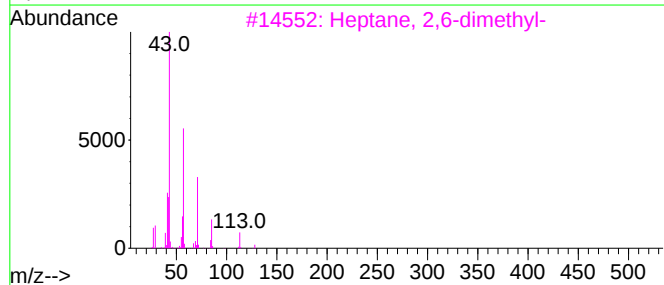
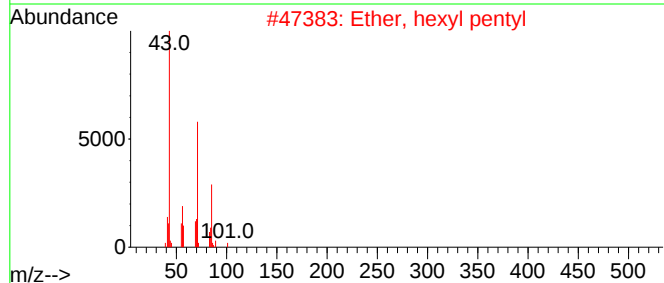
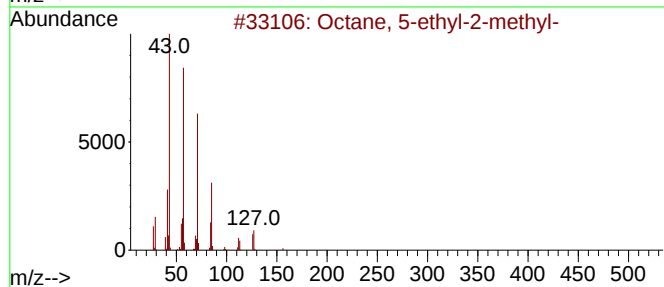
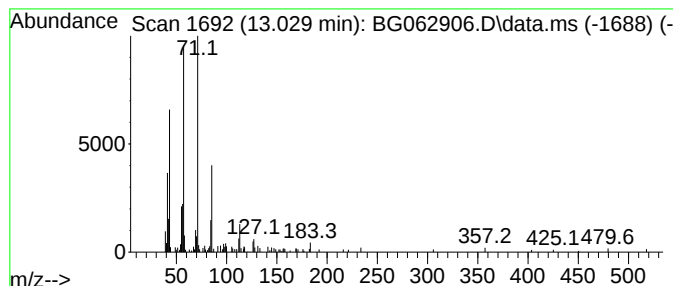
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.029	3.01 ng/ul	119832	Acenaphthene-d10			14.486
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octane, 5-ethyl-2-methyl-		156	C11H24	062016-18-6	78
2	Ether, hexyl pentyl		172	C11H24O	032357-83-8	72
3	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	62
4	Dodecane, 2,6,11-trimethyl-		212	C15H32	031295-56-4	59
5	Octane, 2-methyl-		128	C9H20	003221-61-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

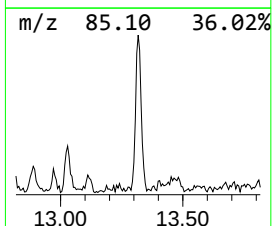
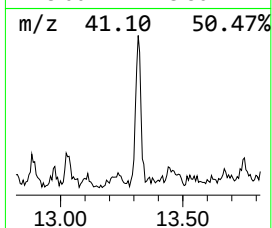
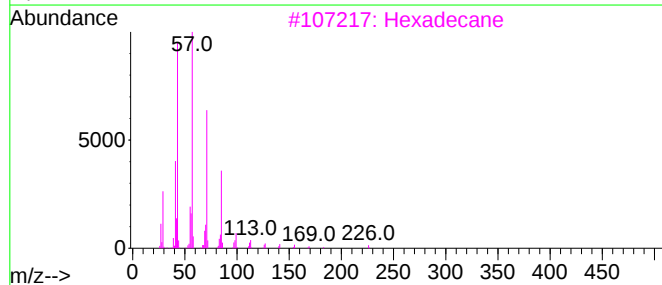
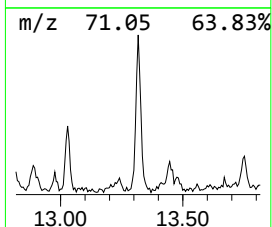
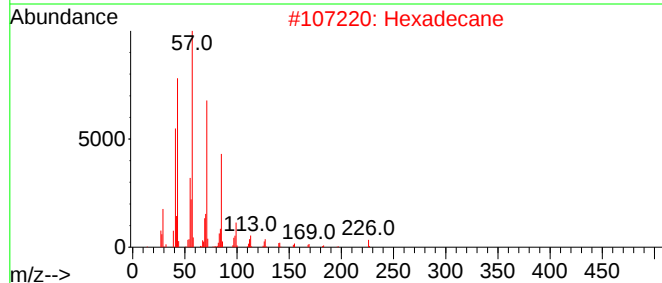
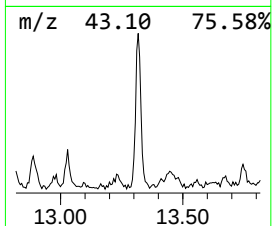
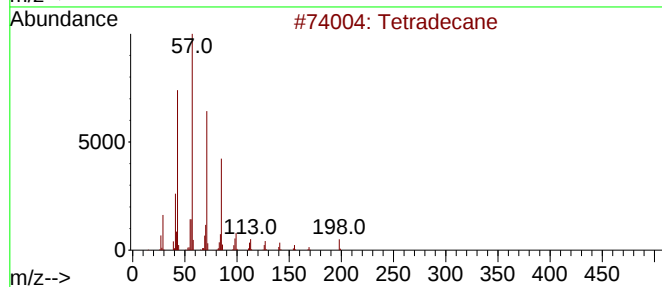
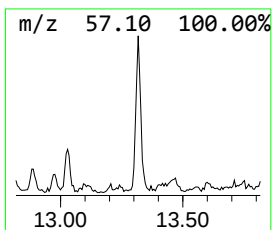
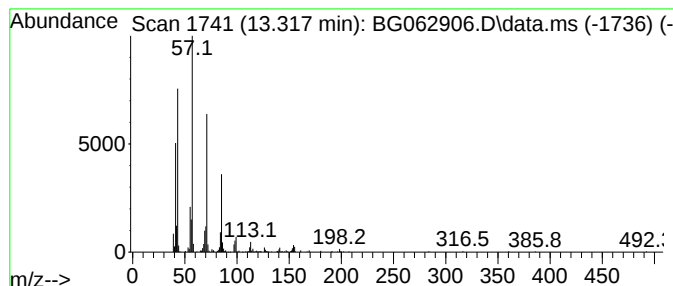
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.317	9.90 ng/ul	394087	Acenaphthene-d10		14.486	
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	Hexadecane		226	C16H34	000544-76-3	87
4	Tridecane		184	C13H28	000629-50-5	86
5	Dodecane		170	C12H26	000112-40-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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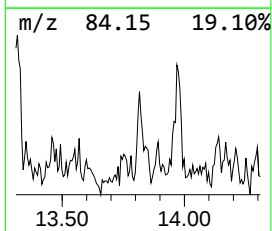
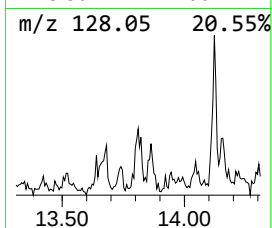
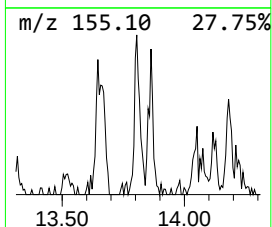
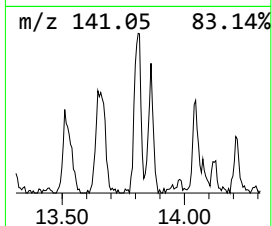
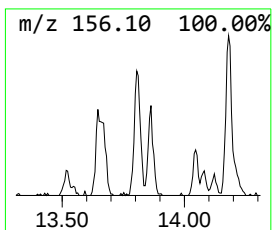
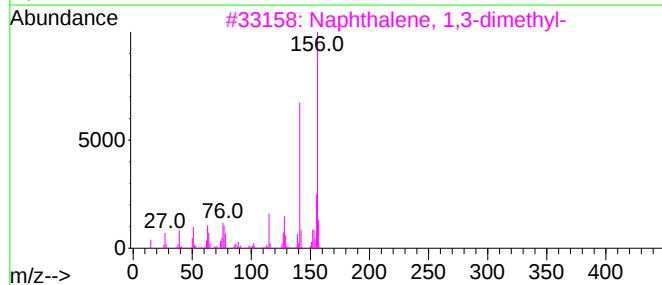
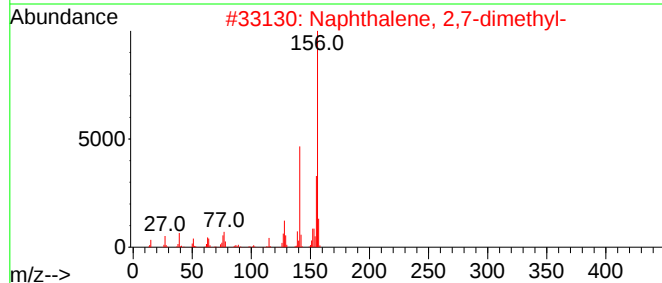
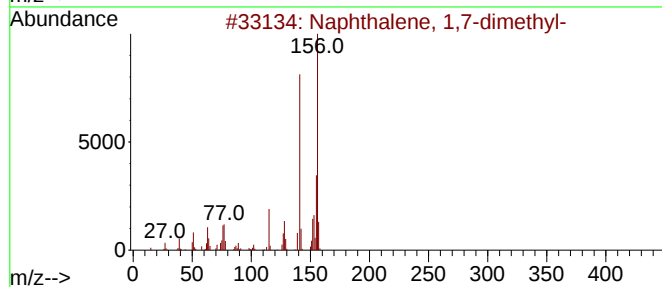
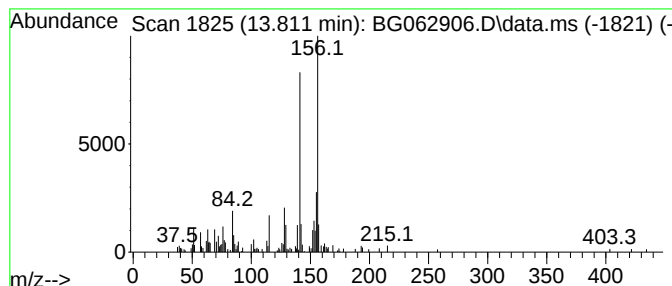
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 Naphthalene, 1,7-dimethyl- Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.811	2.28 ng/ul	90685	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	91
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	91
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	91
4			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	90
5			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

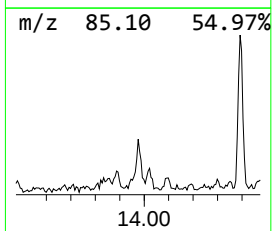
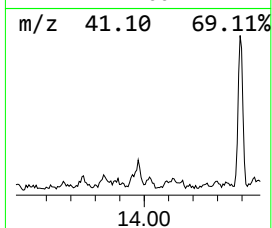
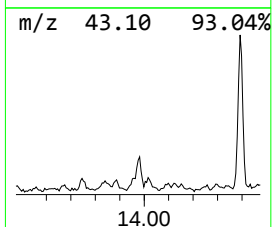
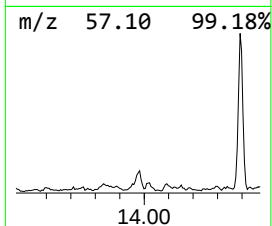
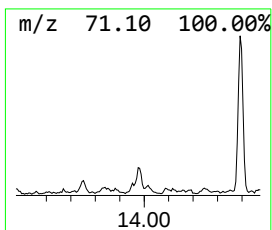
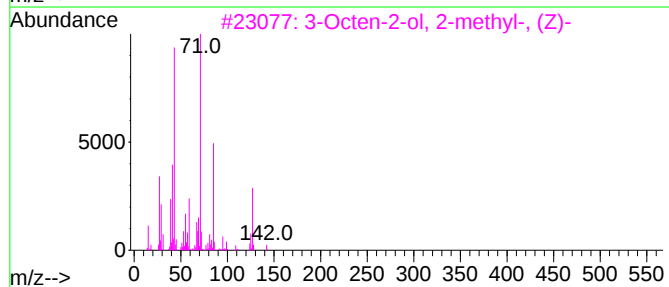
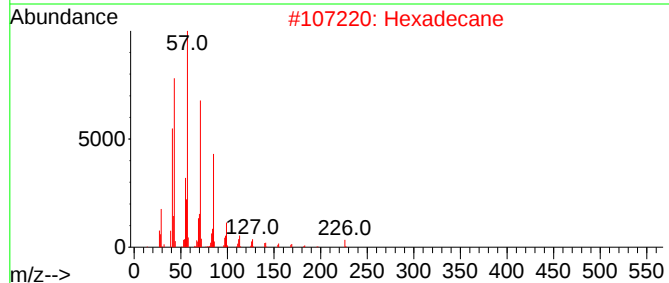
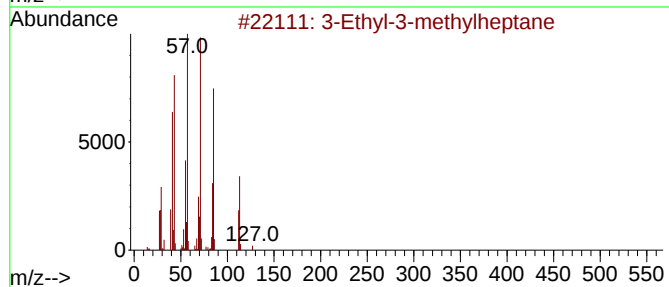
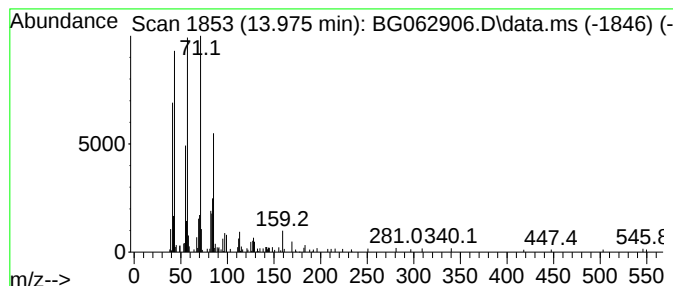
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BNA_G
ClientSampleId :
DDCJ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.975	4.20 ng/ul	167266	Acenaphthene-d10		14.486	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	3-Ethyl-3-methylheptane		142	C10H22	017302-01-1	72
2	Hexadecane		226	C16H34	000544-76-3	59
3	3-Octen-2-ol, 2-methyl-, (Z)-		142	C9H18O	018521-07-8	53
4	4-Heptanone, 3-methyl-		128	C8H16O	015726-15-5	43
5	Decane, 4-ethyl-		170	C12H26	001636-44-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

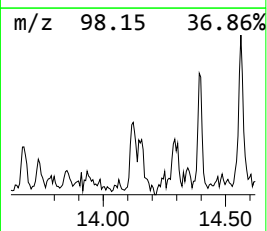
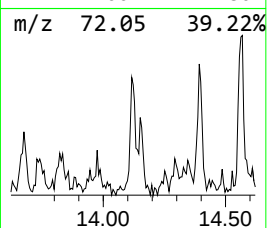
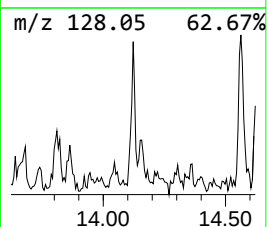
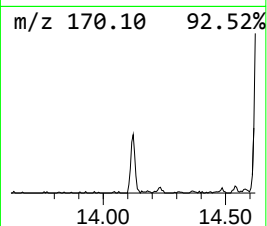
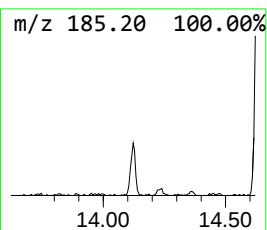
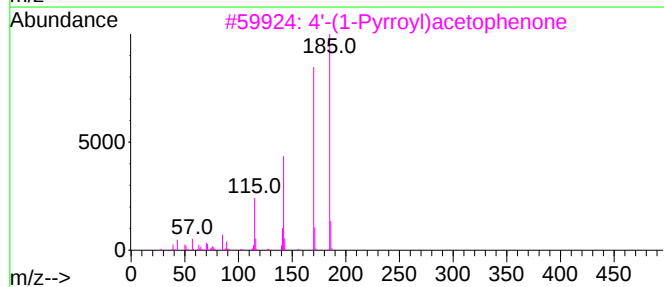
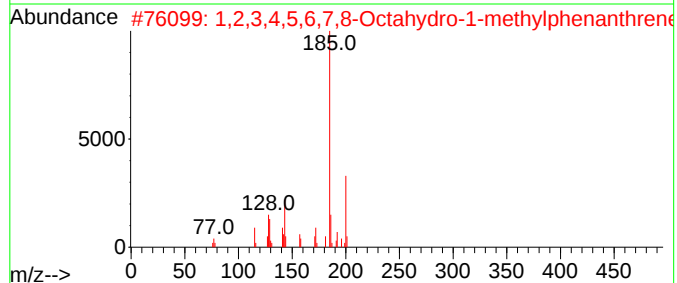
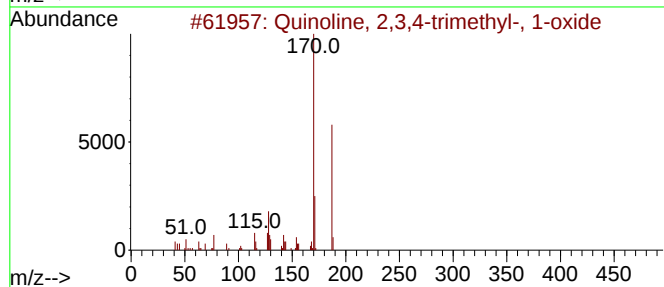
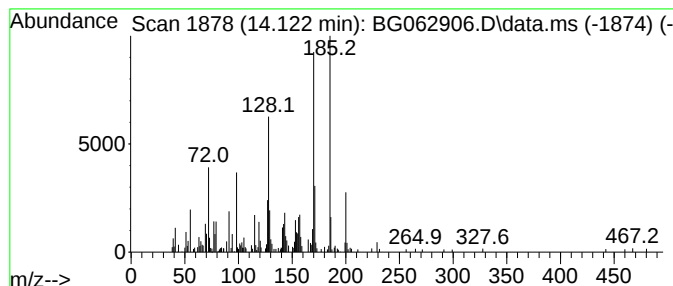
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 Quinoline, 2,3,4-trimethyl-... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.122	3.86 ng/ul	153474	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Quinoline, 2,3,4-trimethyl-, 1-o...	187	C12H13NO	014300-13-1	50
2			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	46
3			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	43
4			3-Hydroxydiphenylamine	185	C12H11NO	000101-18-8	35
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

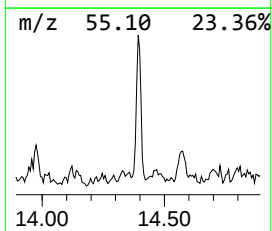
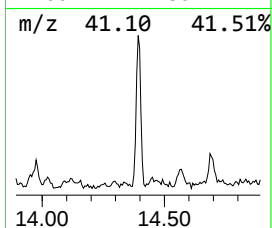
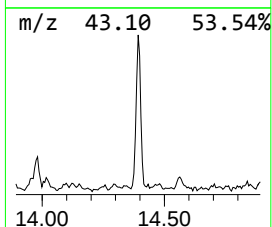
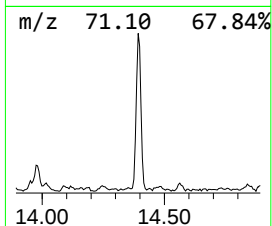
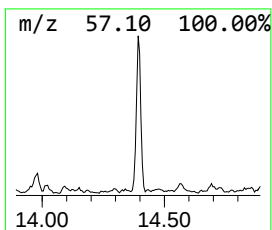
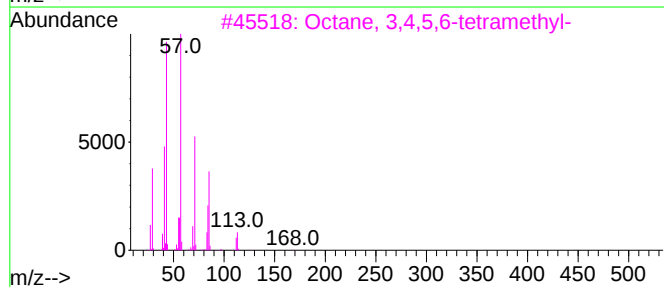
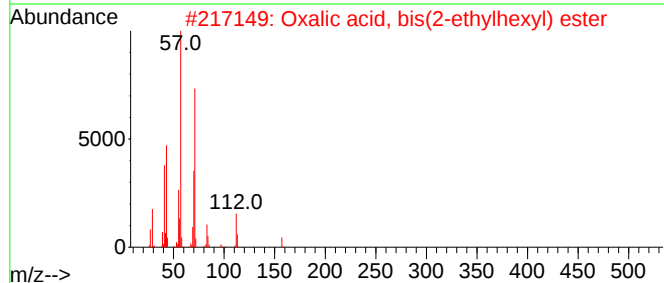
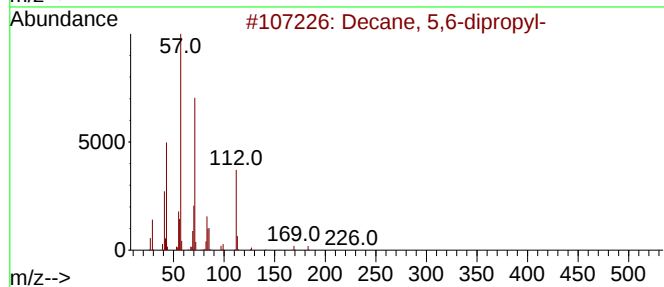
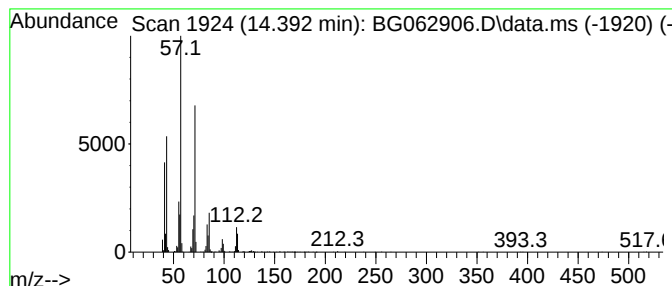
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.392	19.85 ng/ul	789824	Acenaphthene-d10	14.486

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	64
2		Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	64
3		Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	59
4		Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	59
5		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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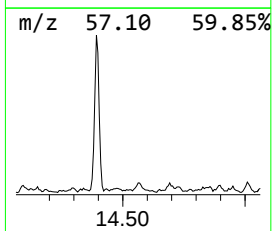
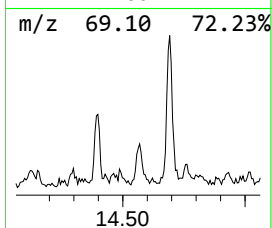
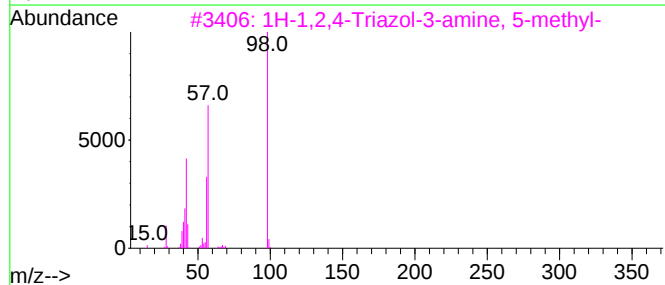
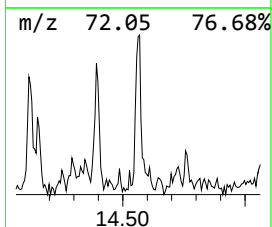
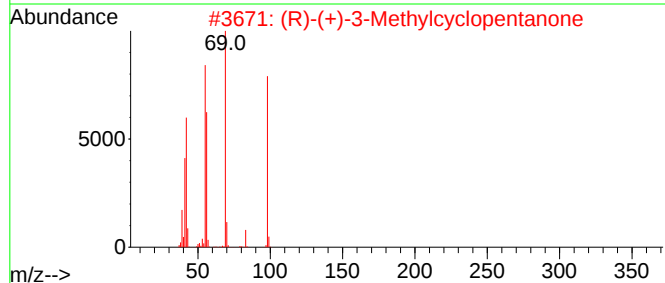
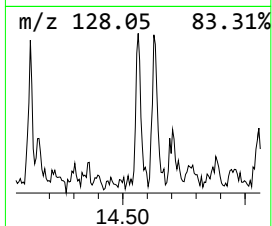
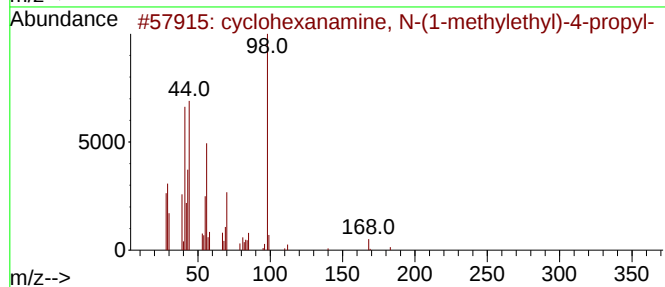
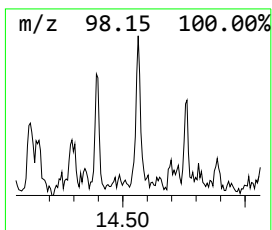
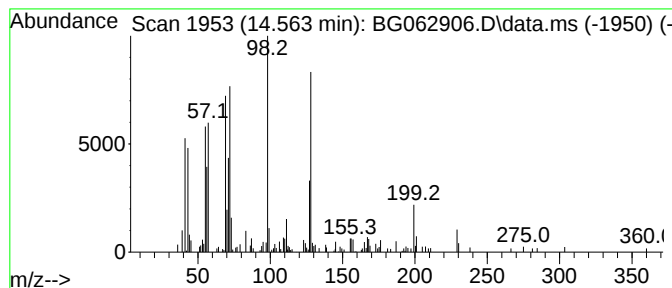
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 unknown-04 Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.563	4.21 ng/ul	167492	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cyclohexanamine, N-(1-methylethy...	183	C12H25N	1000404-90-6	14
2			(R)-(+)-3-Methylcyclopentanone	98	C6H10O	006672-30-6	14
3			1H-1,2,4-Triazol-3-amine, 5-methyl-	98	C3H6N4	004923-01-7	14
4			1,3-Cyclopentanedione, 4-hydroxy...	128	C6H8O3	004800-04-8	14
5			Glycine, N-acetyl-, ethyl ester	145	C6H11NO3	001906-82-7	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

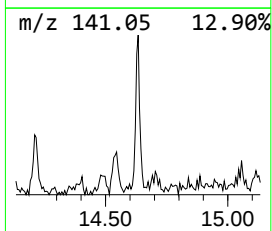
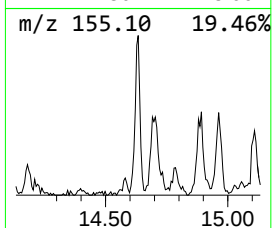
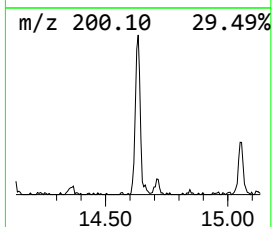
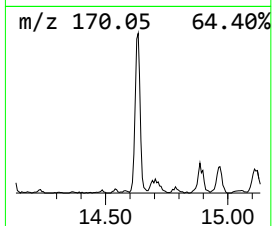
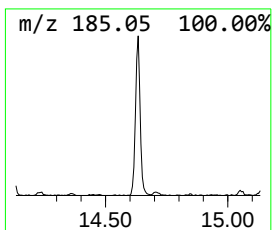
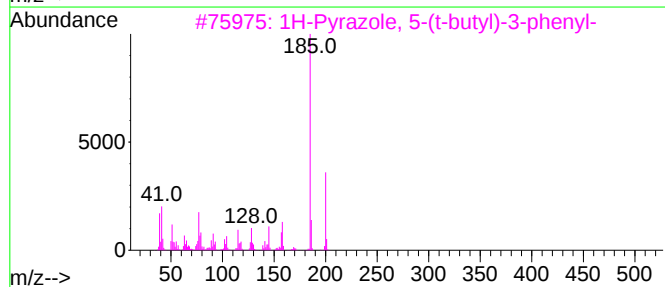
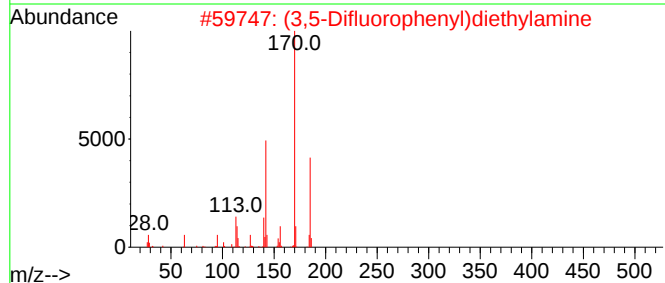
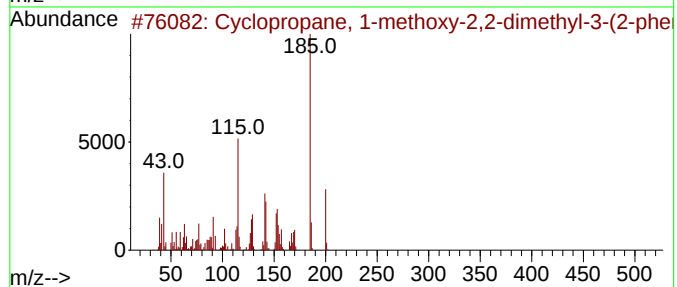
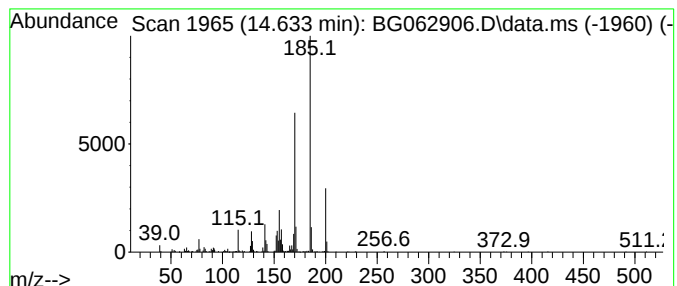
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 Cyclopropane, 1-methoxy-2,2... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.633	11.97 ng/ul	476254	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	64
2			(3,5-Difluorophenyl)diethylamine	185	C10H13F2N	1000306-62-9	52
3			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	49
4			Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	46
5			6-Fluoro-2,4-pyrimidinediamine, TMS	200	C7H13FN4Si	1000472-66-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

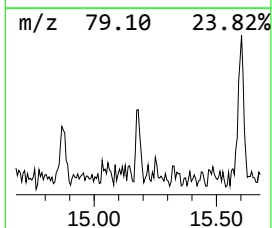
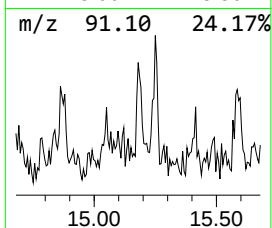
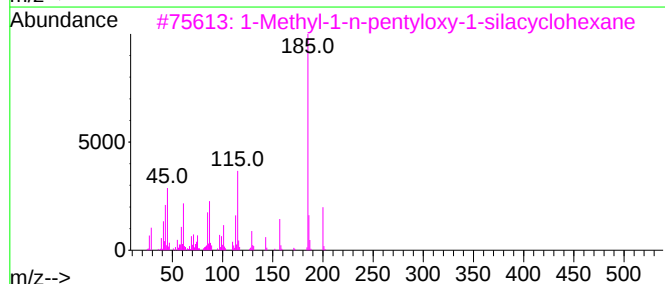
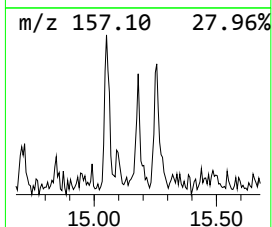
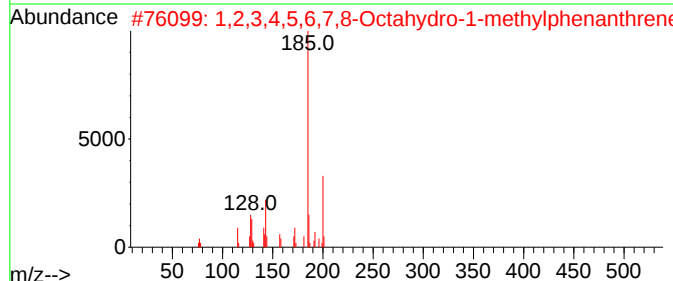
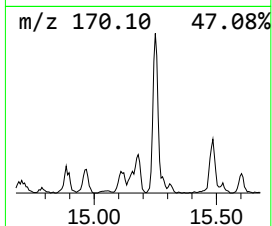
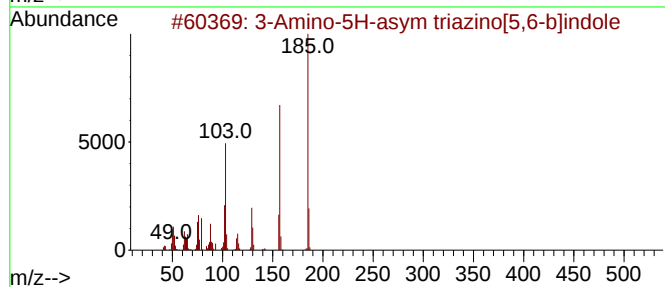
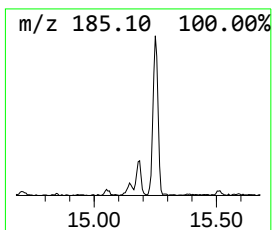
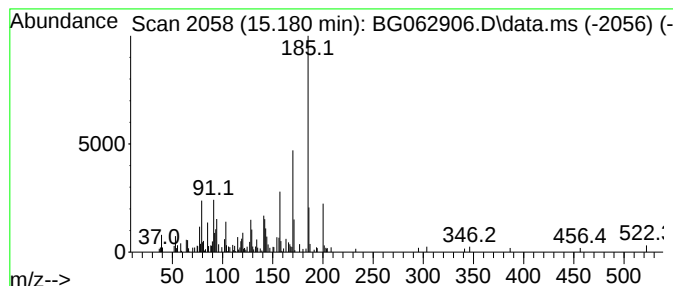
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 3-Amino-5H-asym triazino[5,... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.180	2.77 ng/ul	110283	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Amino-5H-asym triazino[5,6-b]i...	185	C9H7N5	036047-75-3	53
2			1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	49
3			1-Methyl-1-n-pentyloxy-1-silacyc...	200	C11H24OSi	232270-61-0	47
4			Silane, trimethyl-2-naphthalenyl-	200	C13H16Si	018052-85-2	47
5			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

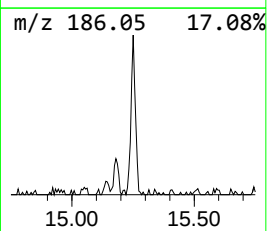
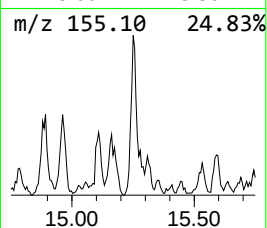
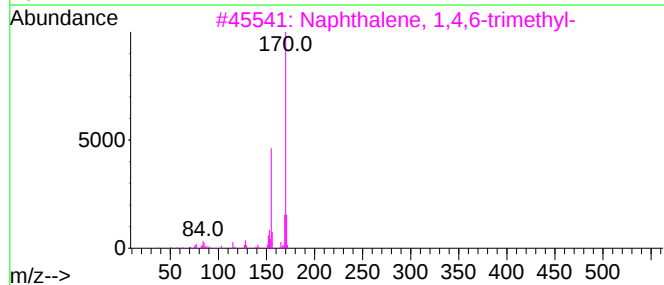
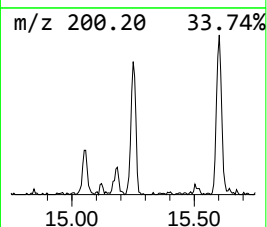
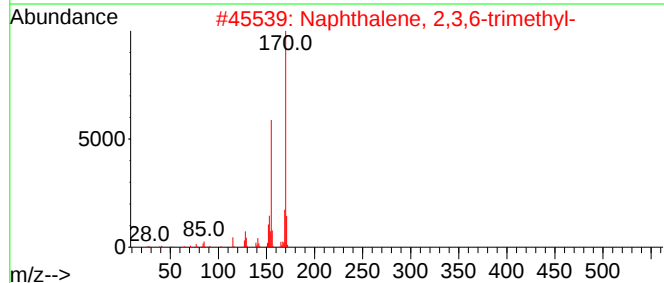
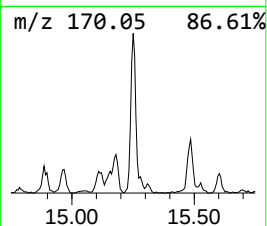
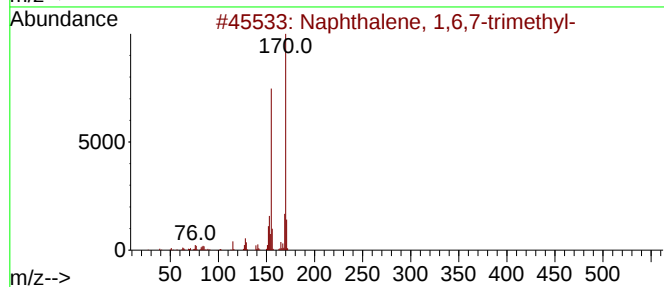
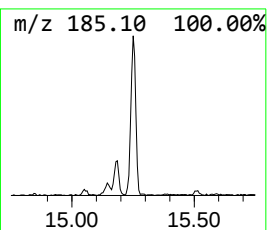
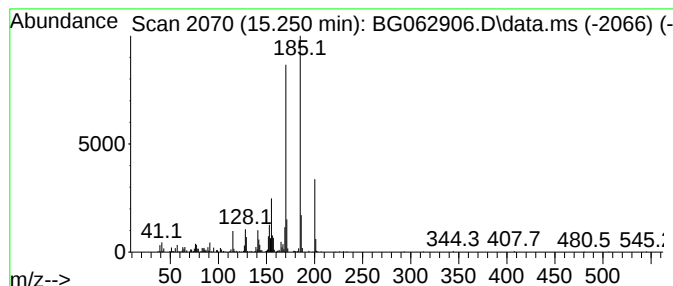
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Naphthalene, 1,6,7-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.250	12.11 ng/ul	481861	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	50
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
4			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	50
5			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

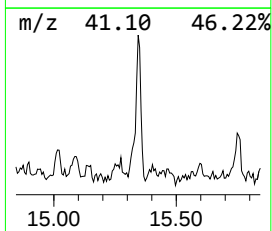
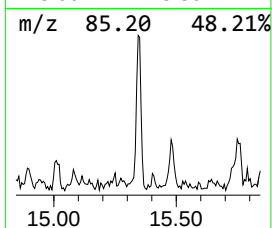
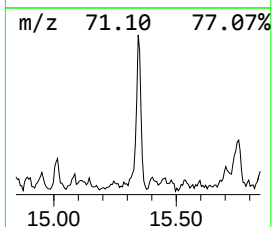
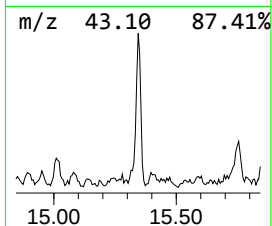
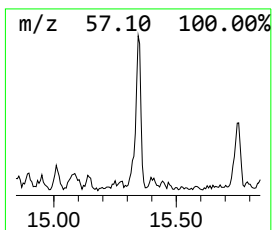
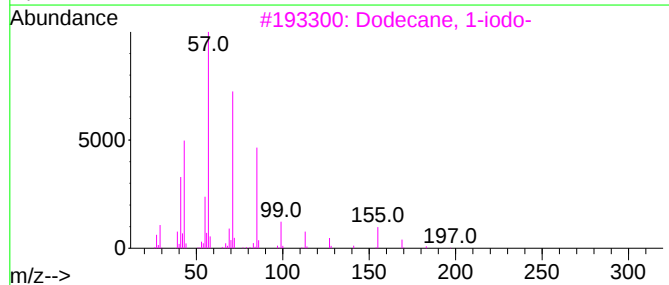
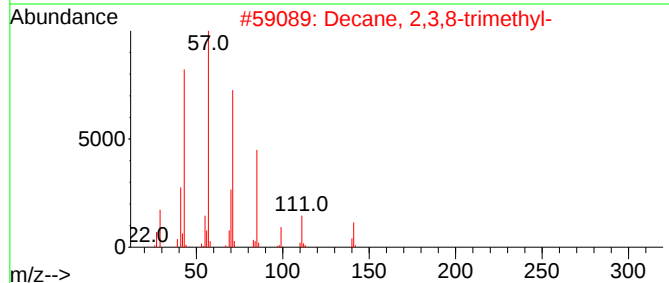
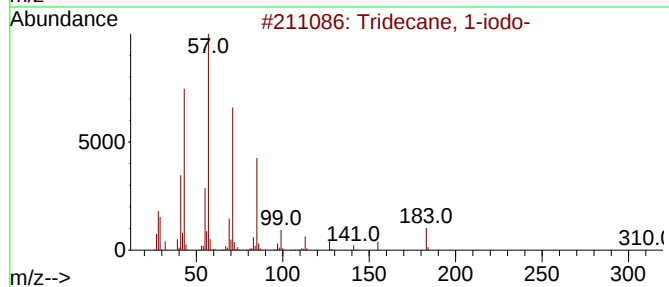
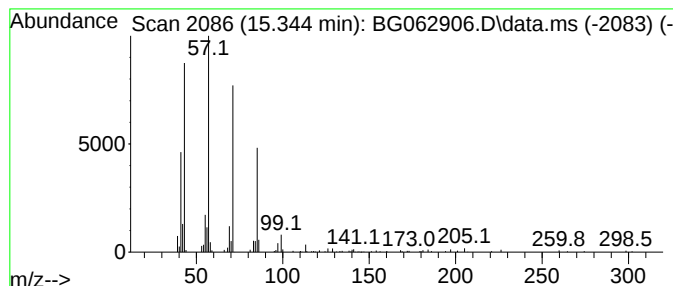
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 Tridecane, 1-iodo- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.344	9.19 ng/ul	365899	Acenaphthene-d10	14.486

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	83
2		Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	78
3		Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
4		Heptacosane	380	C27H56	000593-49-7	72
5		Nonane, 1-iodo-	254	C9H19I	004282-42-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

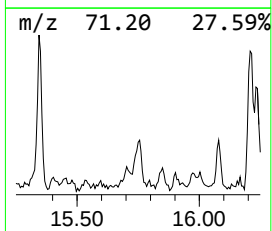
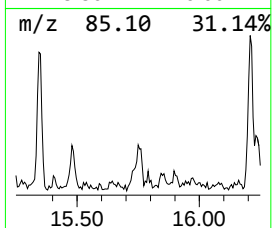
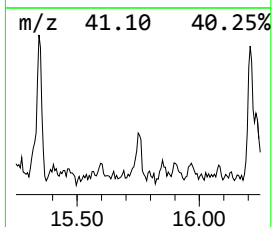
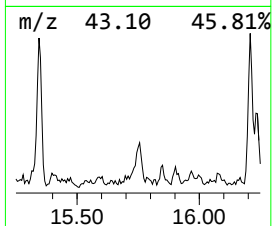
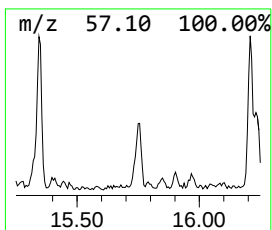
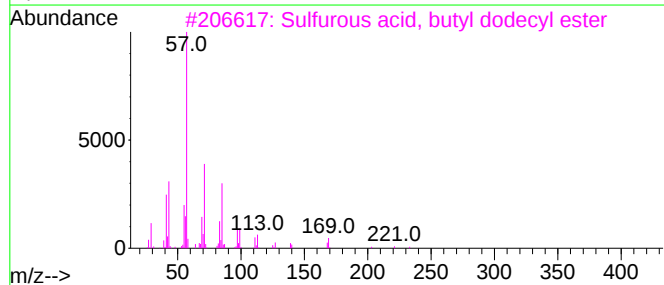
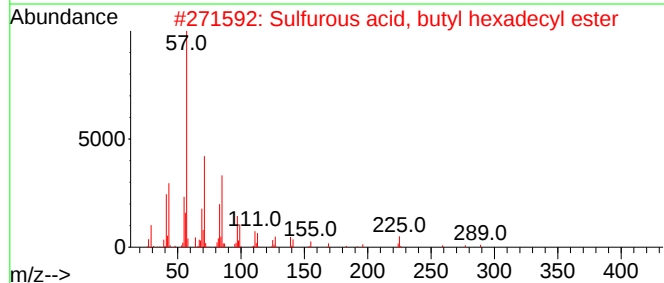
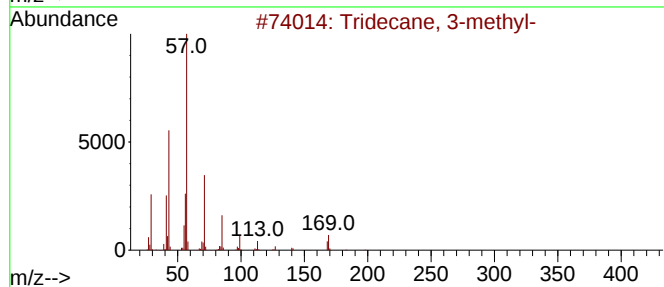
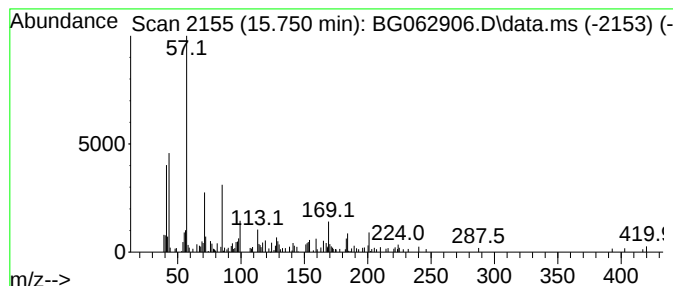
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.749	4.26 ng/ul	169514	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 3-methyl-	198	C14H30	006418-41-3	35
2			Sulfurous acid, butyl hexadecyl ...	362	C20H42O3S	1000309-18-3	27
3			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	25
4			3,5-Dimethyldodecane	198	C14H30	107770-99-0	25
5			3,4-di-n-Butoxy-3-cyclobutene-1,...	226	C12H18O4	002892-62-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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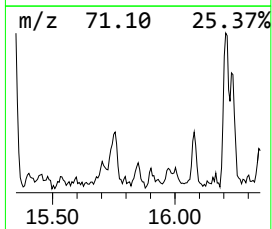
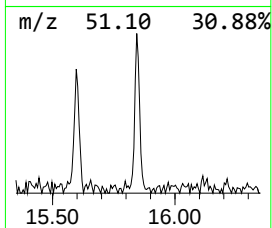
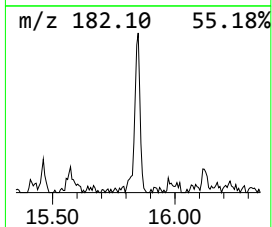
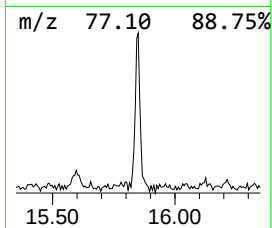
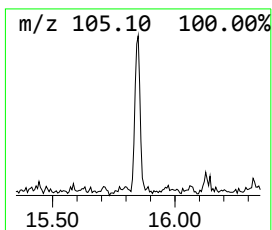
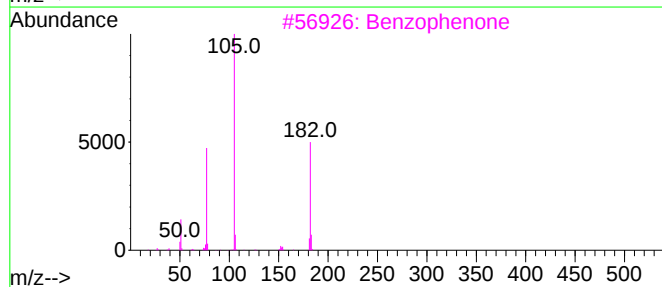
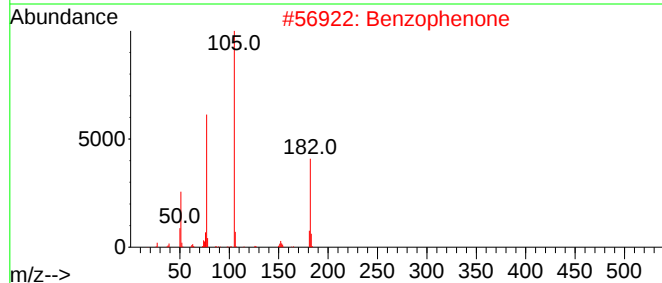
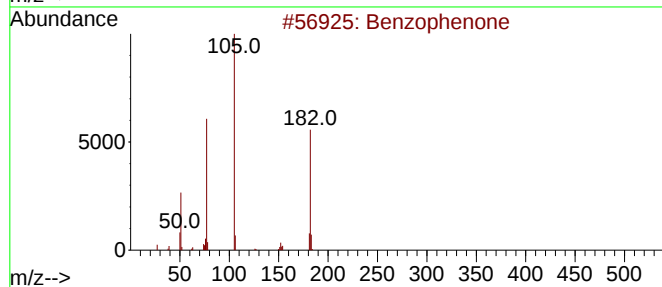
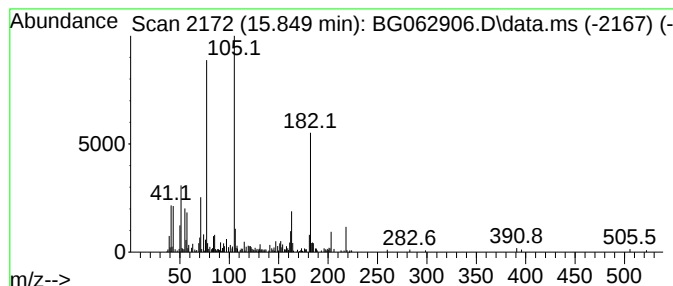
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Benzophenone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.849	5.49 ng/ul	218442	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	70
2			Benzophenone	182	C13H10O	000119-61-9	62
3			Benzophenone	182	C13H10O	000119-61-9	53
4			1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	43
5			1,2-Propanedione, 1-phenyl-, 2-o...	163	C9H9NO2	000119-51-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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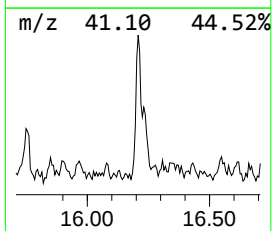
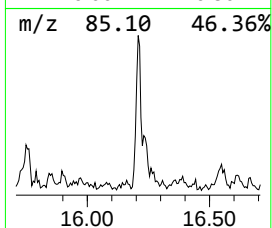
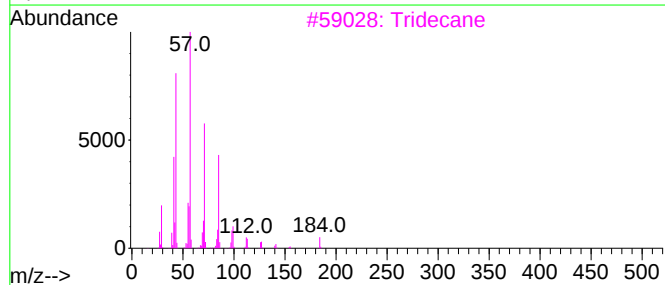
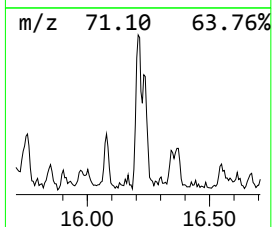
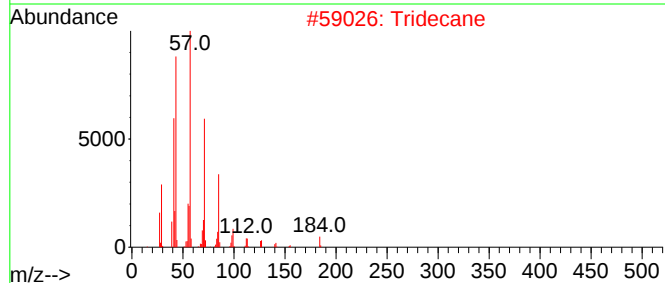
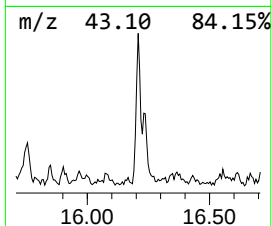
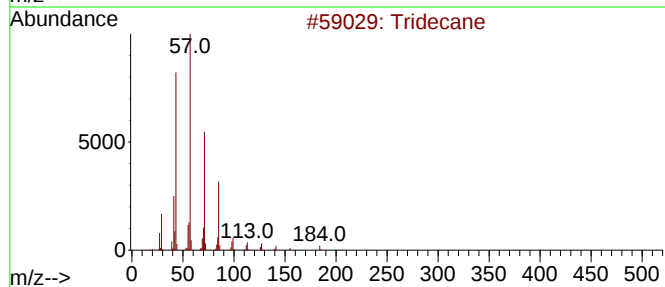
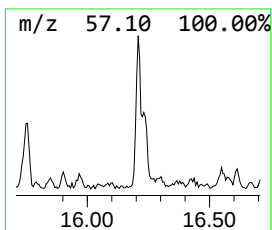
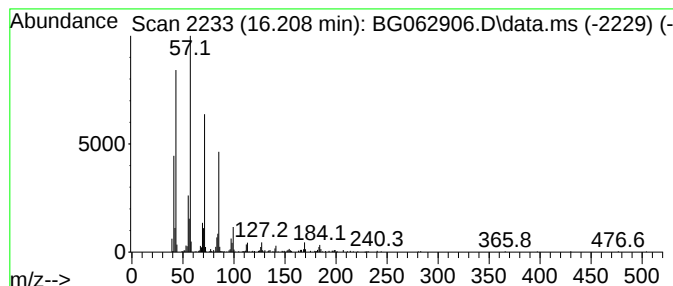
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.208	8.08 ng/ul	416815	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	96
2			Tridecane	184	C13H28	000629-50-5	92
3			Tridecane	184	C13H28	000629-50-5	91
4			Eicosane, 10-methyl-	296	C21H44	054833-23-7	86
5			Eicosane	282	C20H42	000112-95-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

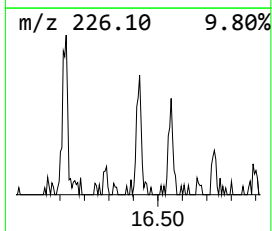
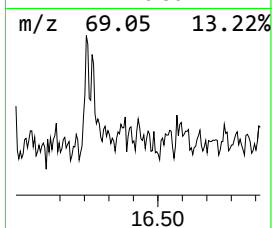
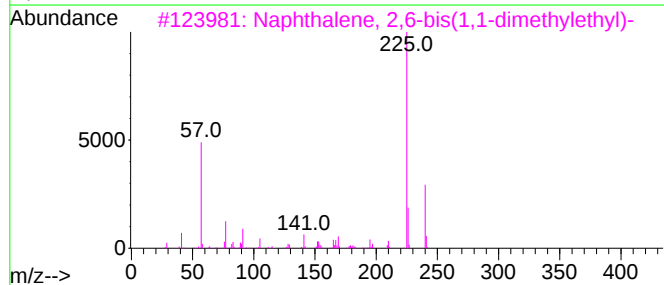
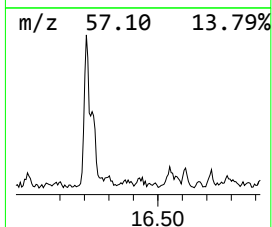
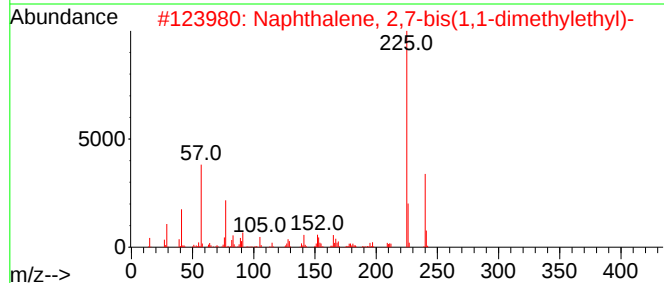
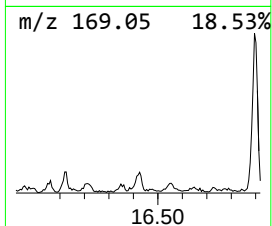
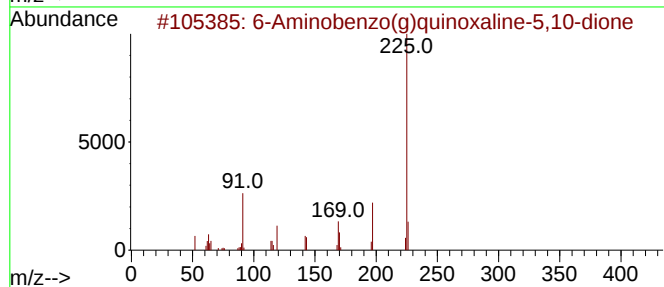
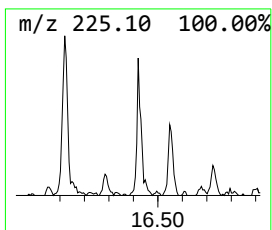
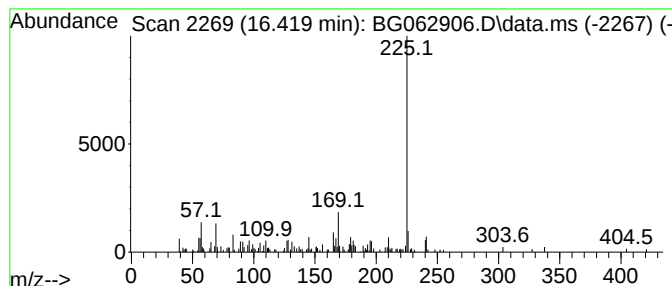
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 42 6-Aminobenzo(g)quinoxaline-... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.419	2.30 ng/ul	118598	Phenanthrene-d10		17.242	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	6-Aminobenzo(g)quinoxaline-5,10-...		225	C12H7N3O2	072225-24-2	53
2	Naphthalene, 2,7-bis(1,1-dimethy...		240	C18H24	010275-58-8	50
3	Naphthalene, 2,6-bis(1,1-dimethy...		240	C18H24	003905-64-4	45
4	N-Methylthioacridone		225	C14H11NS	017435-19-7	38
5	Urea, N-methoxy-N'-(4'-methoxy[1...		272	C15H16N2O3	056667-14-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

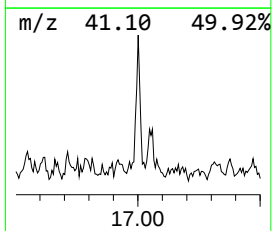
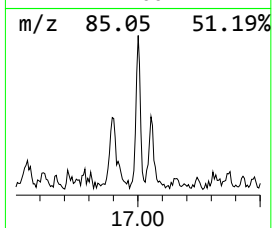
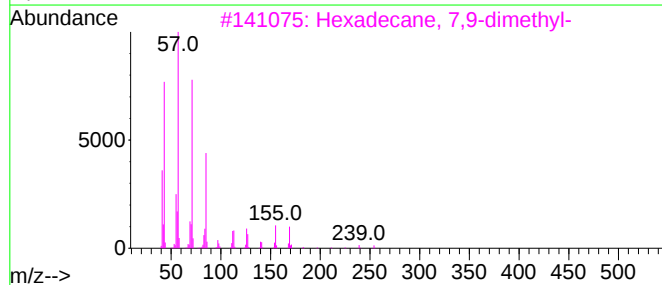
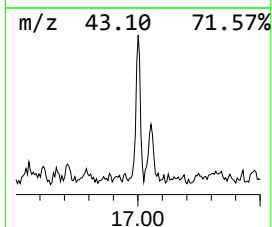
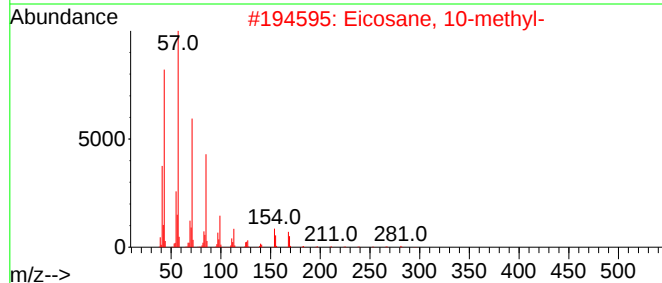
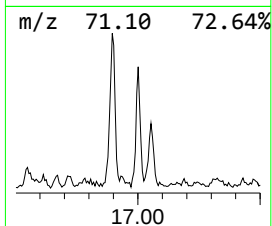
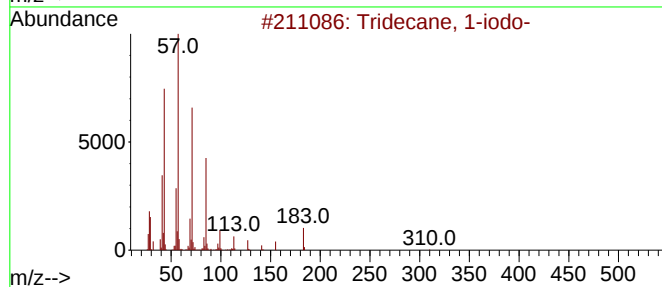
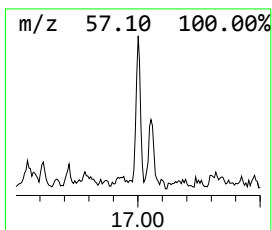
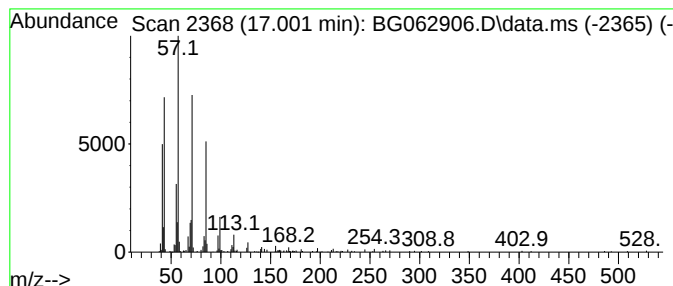
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.001	5.21 ng/ul	268820	Phenanthrene-d10		17.242	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane, 1-iodo-		310	C13H27I	035599-77-0	80
2	Eicosane, 10-methyl-		296	C21H44	054833-23-7	80
3	Hexadecane, 7,9-dimethyl-		254	C18H38	021164-95-4	72
4	Tetracosane		338	C24H50	000646-31-1	72
5	Eicosane		282	C20H42	000112-95-8	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

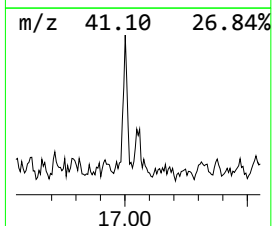
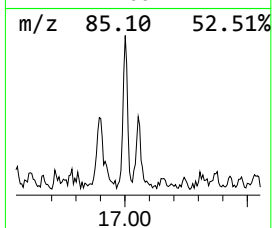
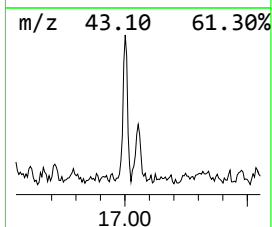
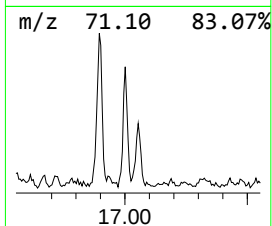
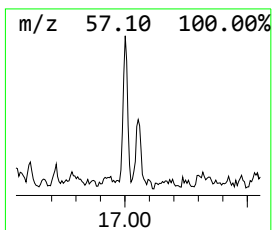
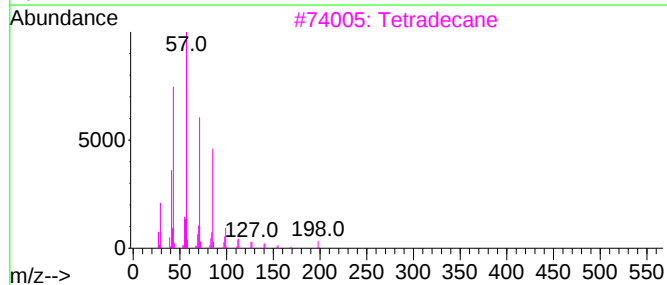
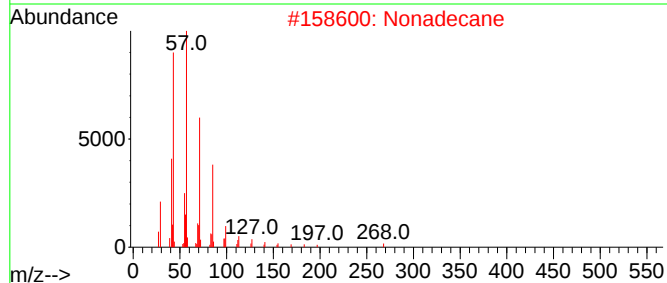
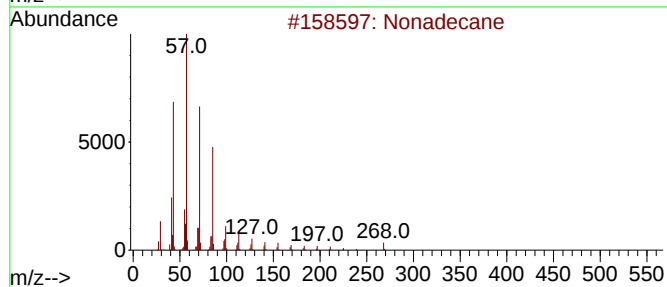
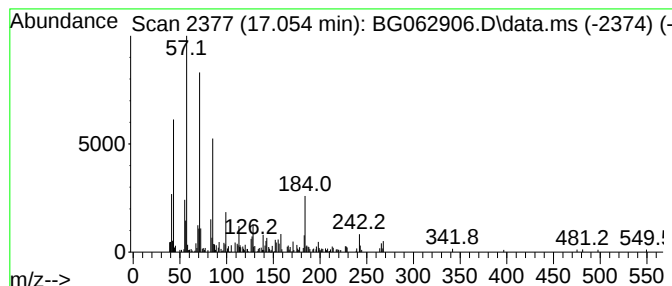
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.054	2.97 ng/ul	153407	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	89
2			Nonadecane	268	C19H40	000629-92-5	68
3			Tetradecane	198	C14H30	000629-59-4	64
4			Sulfurous acid, butyl nonyl ester	264	C13H28O3S	1000309-17-6	58
5			Tetracosane, 1-iodo-	464	C24H49I	1000406-32-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

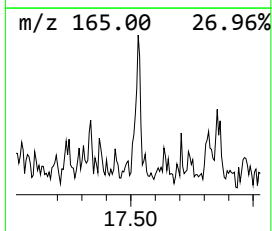
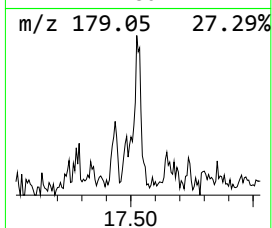
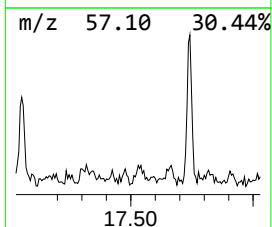
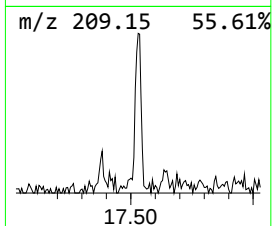
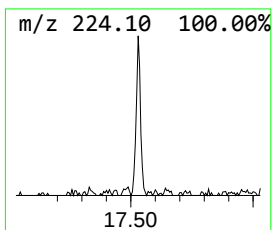
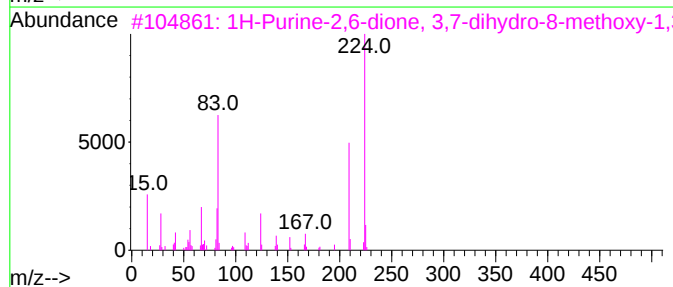
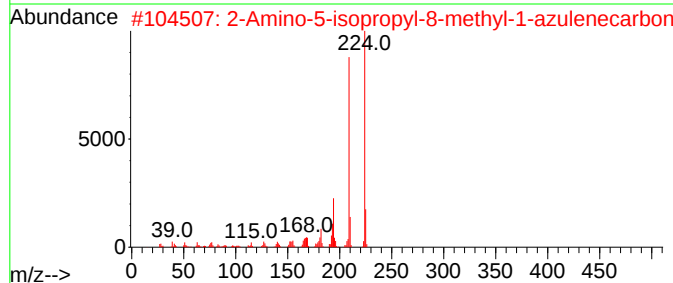
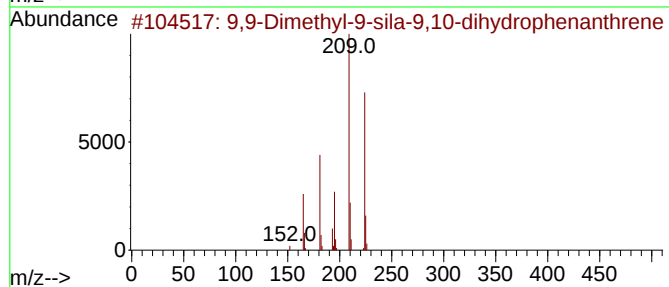
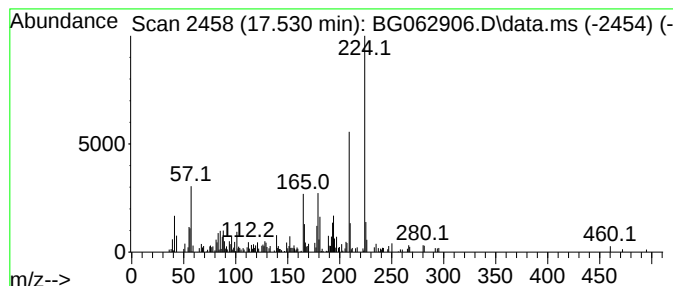
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 9,9-Dimethyl-9-sila-9,10-di... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD			R.T.
17.530	3.96 ng/ul	204535	Phenanthrene-d10			17.242
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,9-Dimethyl-9-sila-9,10-dihydro...	224	C15H16Si		187328-55-8	76
2	2-Amino-5-isopropyl-8-methyl-1-a...	224	C15H16N2		093946-48-6	58
3	1H-Purine-2,6-dione, 3,7-dihydro...	224	C9H12N4O3		000569-34-6	50
4	2-(4-Methoxy-phenyl)-1H-benzoimi...	224	C14H12N2O		1000318-21-9	43
5	Cyclopenta[a]inden-4-one, 6-merc...	224	C9H8N2O52		1000311-03-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

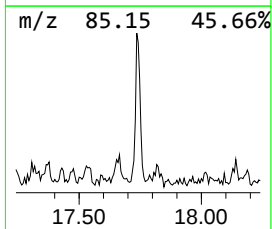
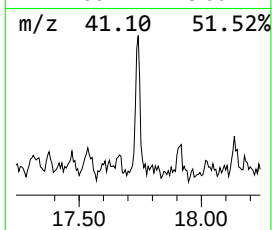
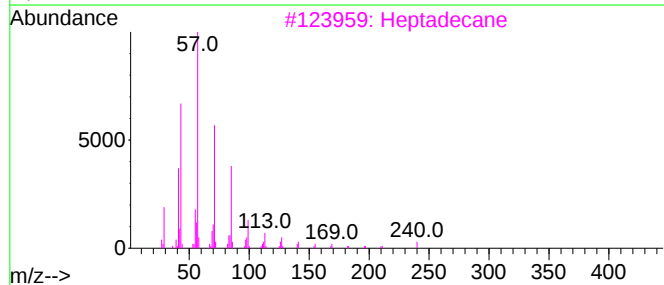
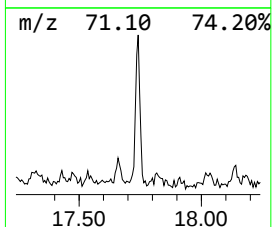
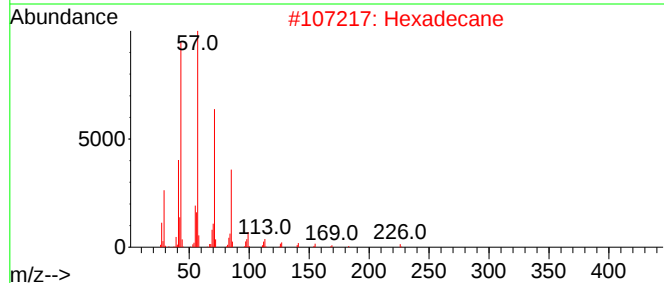
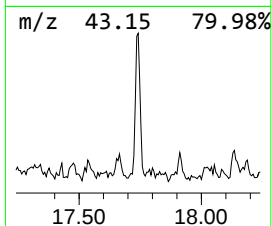
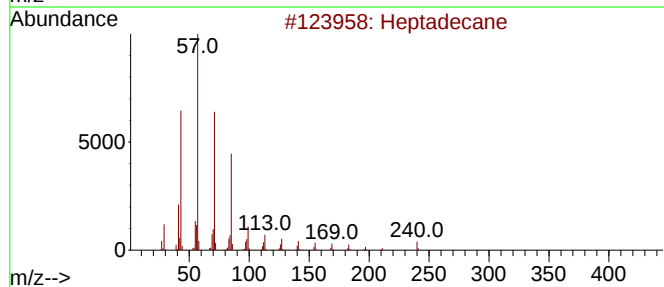
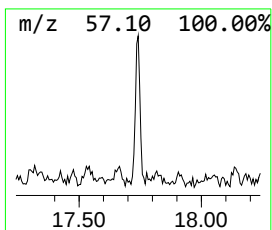
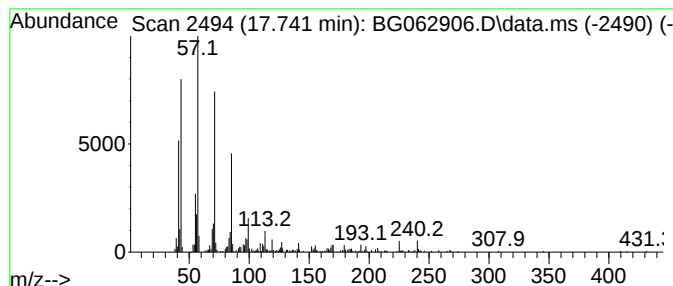
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.741	5.45 ng/ul	281154	Phenanthrene-d10	17.242

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		Heptadecane	240	C17H36	000629-78-7	94
4		Heptadecane	240	C17H36	000629-78-7	94
5		Heptadecane	240	C17H36	000629-78-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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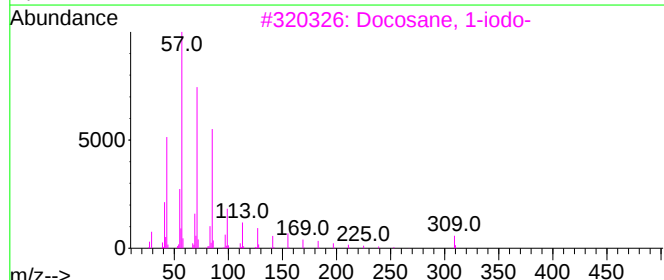
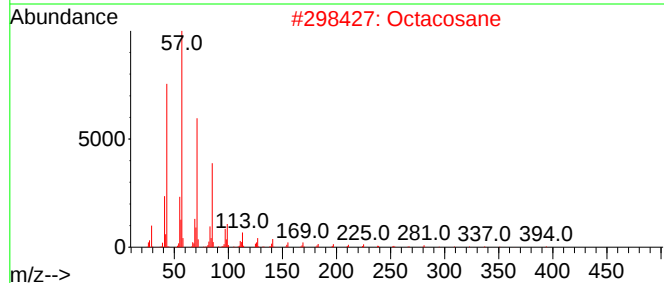
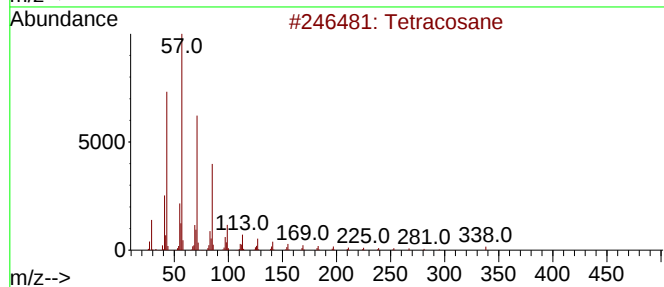
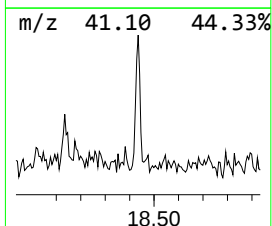
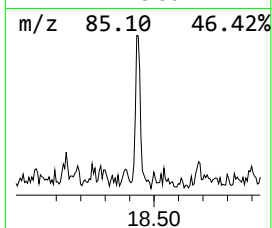
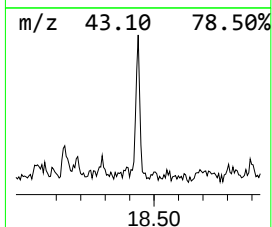
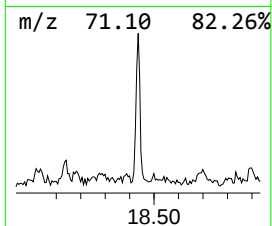
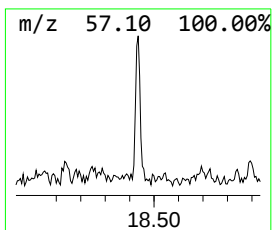
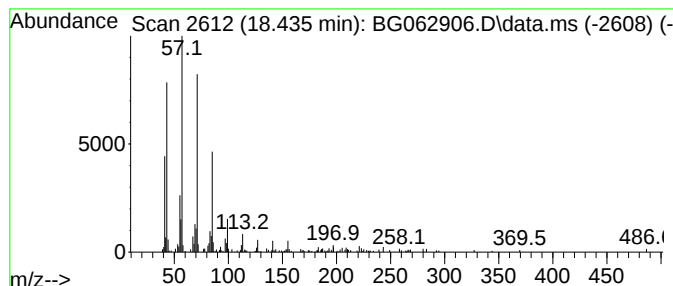
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.435	3.64 ng/ul	187730	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	90
2			Octacosane	394	C28H58	000630-02-4	90
3			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87
4			Heneicosane	296	C21H44	000629-94-7	87
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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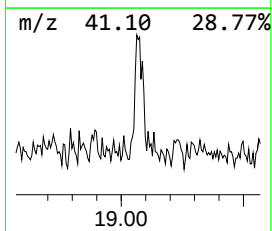
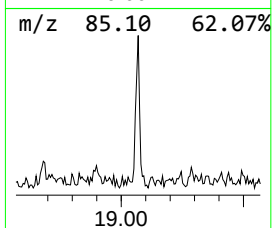
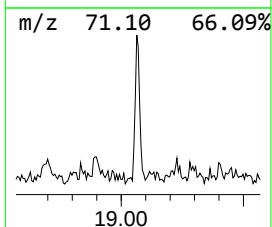
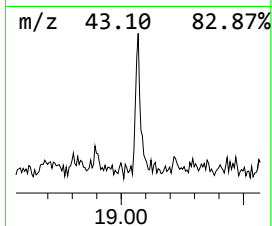
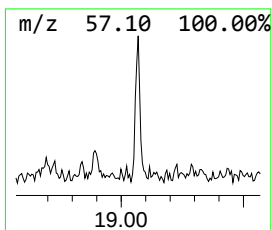
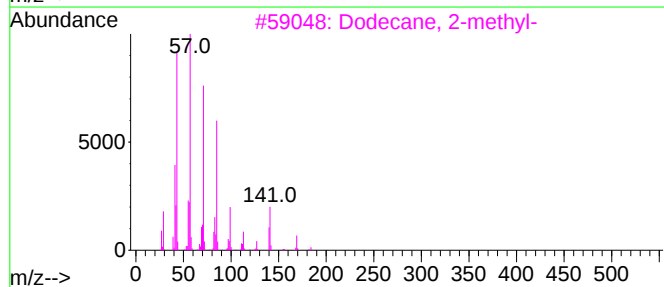
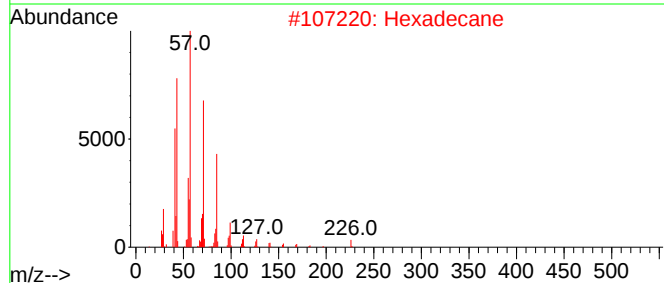
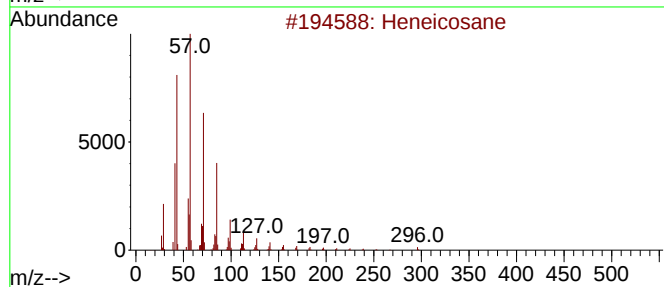
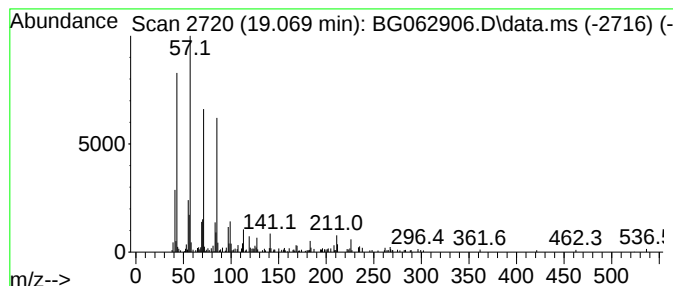
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.069	3.26 ng/ul	167948	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	94
2			Hexadecane	226	C16H34	000544-76-3	93
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	93
4			Heneicosane	296	C21H44	000629-94-7	90
5			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

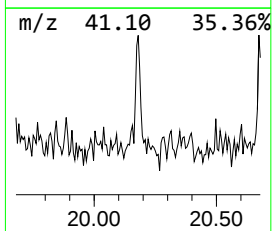
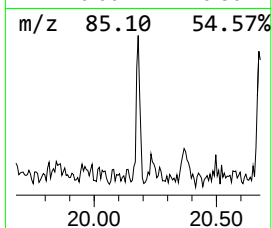
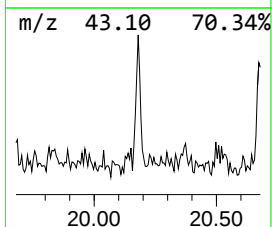
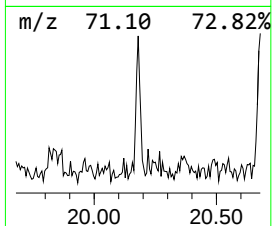
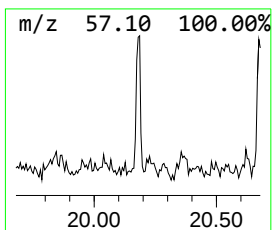
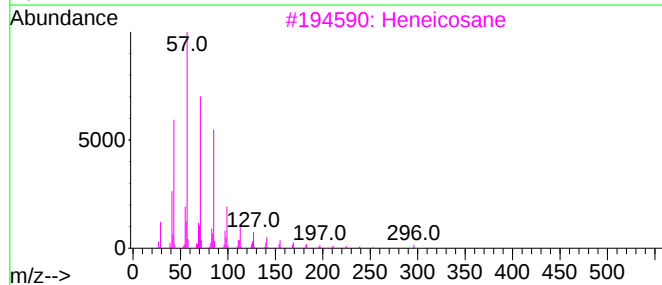
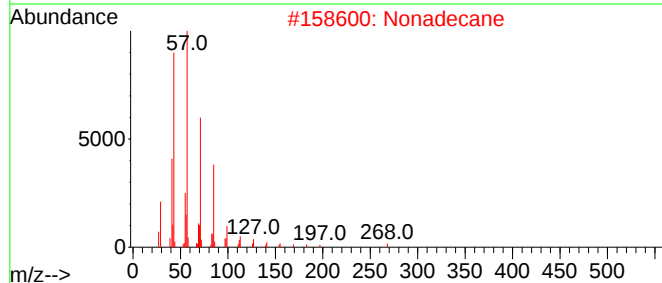
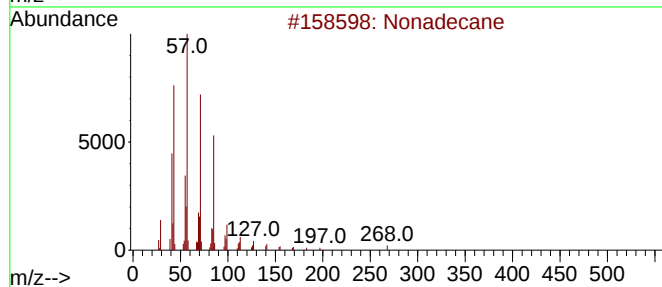
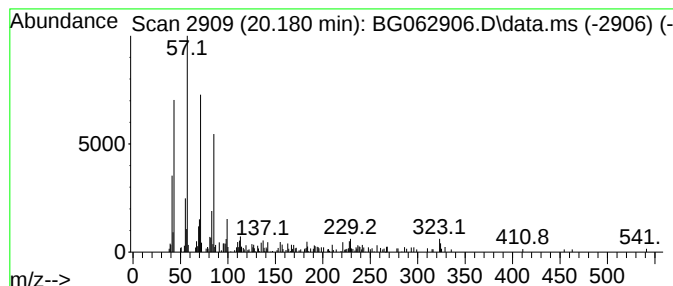
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.180	2.55 ng/ul	137404	Chrysene-d12	21.490

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	78
2		Nonadecane	268	C19H40	000629-92-5	76
3		Heneicosane	296	C21H44	000629-94-7	76
4		Tricosane	324	C23H48	000638-67-5	72
5		Docosane	310	C22H46	000629-97-0	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062906.D
Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDCJ0ME

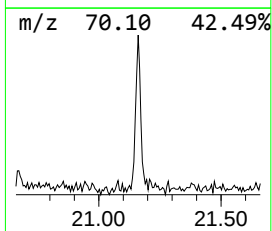
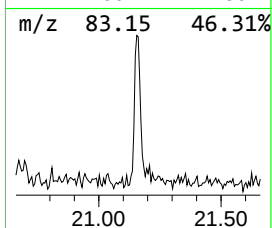
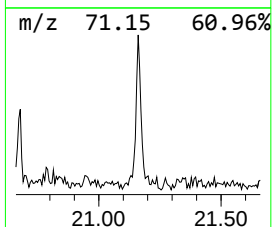
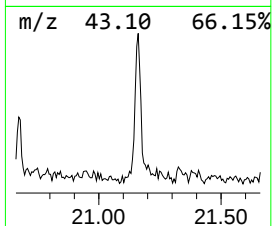
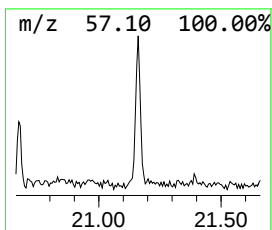
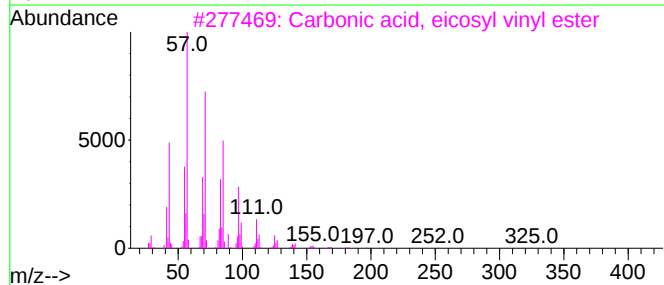
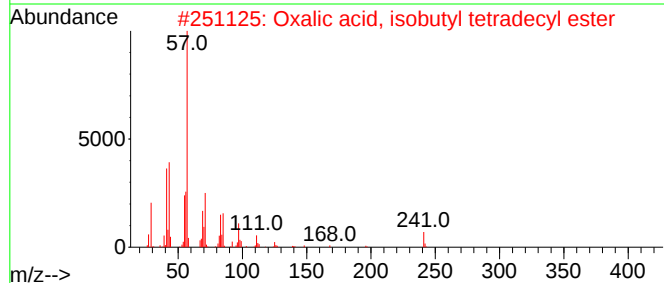
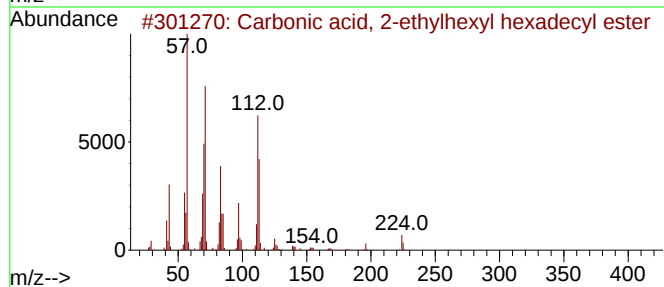
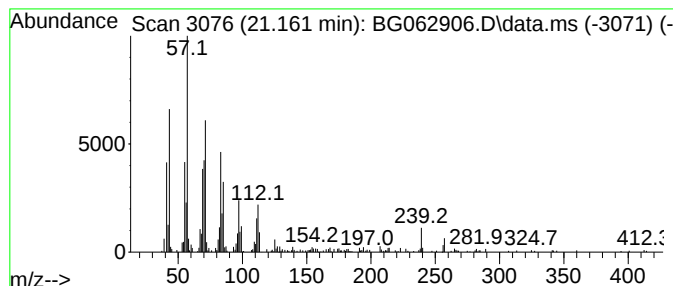
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Quant Title : SVOA CALIBRATION

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

Peak Number 50 Carbonic acid, 2-ethylhexyl... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.161	6.67 ng/ul	360214	Chrysene-d12	21.490

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, 2-ethylhexyl hexa...	398	C25H50O3	1000383-14-3	64
2			Oxalic acid, isobutyl tetradecyl...	342	C20H38O4	1000309-37-9	64
3			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	49
4			Oxalic acid, 2-ethylhexyl isohex...	286	C16H30O4	1000309-38-8	49
5			Eicosyl octyl ether	410	C28H58O	1000406-38-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
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Acq On : 18 Sep 2024 13:11
Operator : JU/RC
Sample : P3878-07ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

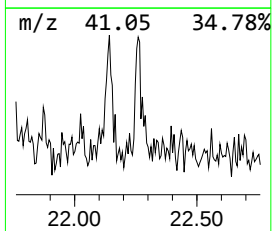
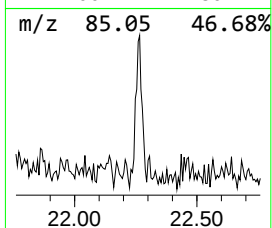
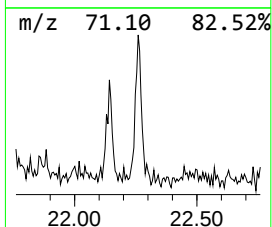
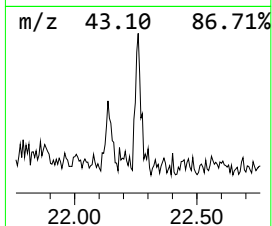
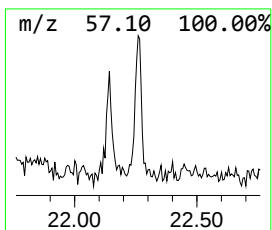
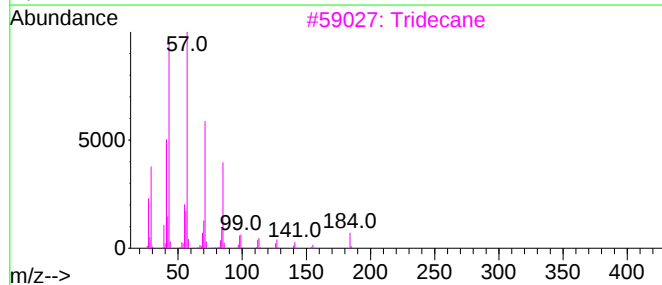
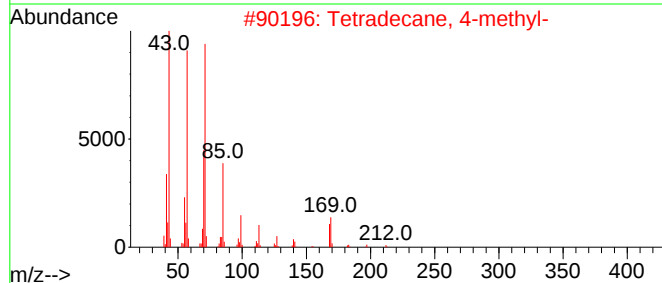
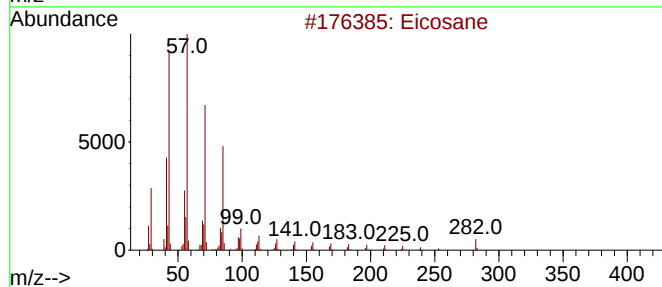
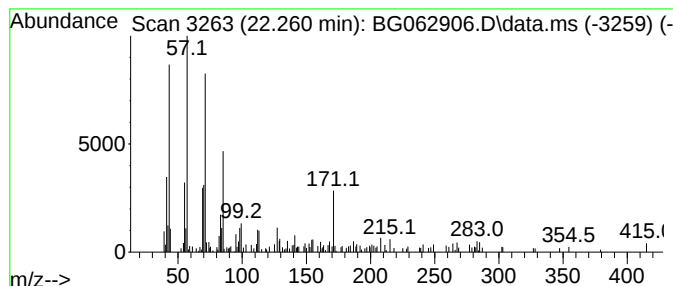
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BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.259	2.89 ng/ul	155774	Chrysene-d12		21.490	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	83
2	Tetradecane, 4-methyl-		212	C15H32	025117-24-2	83
3	Tridecane		184	C13H28	000629-50-5	64
4	Eicosane		282	C20H42	000112-95-8	64
5	Docosane		310	C22H46	000629-97-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062906.D
 Acq On : 18 Sep 2024 13:11
 Operator : JU/RC
 Sample : P3878-07ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDCJ0ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.885	8.5	ng/ul	182289	1	7.859	429077	20.0
(DEL) Alkane: S...	6.425	2.8	ng/ul	60772	1	7.859	429077	20.0
(DEL) Alkane: S...	6.860	4.4	ng/ul	94879	1	7.859	429077	20.0
(DEL) Alkane: S...	6.913	5.3	ng/ul	113592	1	7.859	429077	20.0
Benmoxin	6.954	3.3	ng/ul	70795	1	7.859	429077	20.0
(DEL) Alkane: S...	7.489	29.5	ng/ul	632780	1	7.859	429077	20.0
unknown-01	7.777	2.8	ng/ul	59254	1	7.859	429077	20.0
1-Hexanol, 2-et...	7.918	11.7	ng/ul	251757	1	7.859	429077	20.0
Benzene, 1,2,4-...	7.976	2.7	ng/ul	58290	1	7.859	429077	20.0
1,3-Cyclohexane...	8.288	3.0	ng/ul	64128	1	7.859	429077	20.0
(DEL) Alkane: S...	8.393	5.7	ng/ul	122533	1	7.859	429077	20.0
Sulfurous acid,...	8.446	2.6	ng/ul	56836	1	7.859	429077	20.0
(DEL) Alkane: S...	8.623	4.3	ng/ul	93367	1	7.859	429077	20.0
Benzene, 1,2,3,...	8.805	2.6	ng/ul	54998	1	7.859	429077	20.0
Benzene, 1,2,4,...	8.858	3.4	ng/ul	72354	1	7.859	429077	20.0
(DEL) Alkane: S...	9.087	29.1	ng/ul	624369	1	7.859	429077	20.0
unknown-02	9.204	7.0	ng/ul	150250	1	7.859	429077	20.0
(DEL) Alkane: S...	9.792	2.5	ng/ul	129929	2	10.644	1049530	20.0
Cyanic acid, et...	11.026	8.9	ng/ul	468724	2	10.644	1049530	20.0
unknown-03	11.149	2.9	ng/ul	150568	2	10.644	1049530	20.0
(DEL) Alkane: S...	11.678	4.0	ng/ul	209570	2	10.644	1049530	20.0
Benzene, 4-(2-b...	11.743	7.3	ng/ul	384398	2	10.644	1049530	20.0
2,4-Dipropyl-5-...	11.907	4.6	ng/ul	242336	2	10.644	1049530	20.0
(DEL) Alkane: S...	12.072	7.1	ng/ul	374688	2	10.644	1049530	20.0
(DEL) Alkane: S...	12.107	4.6	ng/ul	241658	2	10.644	1049530	20.0
(DEL) Alkane: S...	12.888	2.3	ng/ul	91168	3	14.486	795871	20.0
(DEL) Alkane: S...	13.029	3.0	ng/ul	119832	3	14.486	795871	20.0
(DEL) Alkane: S...	13.317	9.9	ng/ul	394087	3	14.486	795871	20.0
Naphthalene, 1,...	13.811	2.3	ng/ul	90685	3	14.486	795871	20.0
(DEL) Alkane: S...	13.975	4.2	ng/ul	167266	3	14.486	795871	20.0
Quinoline, 2,3,...	14.122	3.9	ng/ul	153474	3	14.486	795871	20.0
(DEL) Alkane: S...	14.392	19.9	ng/ul	789824	3	14.486	795871	20.0
unknown-04	14.563	4.2	ng/ul	167492	3	14.486	795871	20.0
Cyclopropane, 1...	14.633	12.0	ng/ul	476254	3	14.486	795871	20.0
3-Amino-5H-asym...	15.180	2.8	ng/ul	110283	3	14.486	795871	20.0
Naphthalene, 1,...	15.250	12.1	ng/ul	481861	3	14.486	795871	20.0
Tridecane, 1-iodo-	15.344	9.2	ng/ul	365899	3	14.486	795871	20.0
(DEL) Alkane: S...	15.749	4.3	ng/ul	169514	3	14.486	795871	20.0
Benzophenone	15.849	5.5	ng/ul	218442	3	14.486	795871	20.0
2-Methylidenehy...	16.120	3.6	ng/ul	186783	4	17.242	1031900	20.0
(DEL) Alkane: S...	16.208	8.1	ng/ul	416815	4	17.242	1031900	20.0
6-Aminobenzo(g)...	16.419	2.3	ng/ul	118598	4	17.242	1031900	20.0
(DEL) Alkane: S...	17.001	5.2	ng/ul	268820	4	17.242	1031900	20.0
(DEL) Alkane: S...	17.054	3.0	ng/ul	153407	4	17.242	1031900	20.0
9,9-Dimethyl-9-...	17.530	4.0	ng/ul	204535	4	17.242	1031900	20.0
(DEL) Alkane: S...	17.741	5.5	ng/ul	281154	4	17.242	1031900	20.0
(DEL) Alkane: S...	18.435	3.6	ng/ul	187730	4	17.242	1031900	20.0
(DEL) Alkane: S...	19.069	3.3	ng/ul	167948	4	17.242	1031900	20.0
(DEL) Alkane: S...	20.180	2.5	ng/ul	137404	5	21.490	1079660	20.0
Carbonic acid, ...	21.161	6.7	ng/ul	360214	5	21.490	1079660	20.0
(DEL) Alkane: S...	22.259	2.9	ng/ul	155774	5	21.490	1079660	20.0