

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062969.D
 Acq On : 20 Sep 2024 8:39
 Operator : JU/RC
 Sample : P3898-20ME
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC59ME

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: BG062969.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.741	106	111	121	rBV2	78292	153198	1.15%	0.293%
2	5.885	469	476	485	rVB2	321842	528884	3.98%	1.011%
3	6.355	550	556	563	rBV6	85209	169180	1.27%	0.323%
4	6.426	563	568	572	rBV2	101541	153280	1.15%	0.293%
5	6.508	572	582	584	rBV4	85935	183853	1.38%	0.351%
6	6.855	630	641	646	rBV3	166023	394369	2.97%	0.754%
7	6.913	646	651	654	rBV2	176791	263197	1.98%	0.503%
8	6.949	654	657	662	rVB	186163	283686	2.13%	0.542%
9	7.019	662	669	676	rVB4	349333	720951	5.42%	1.378%
10	7.084	676	680	685	rVB	120443	181416	1.36%	0.347%
11	7.189	693	698	702	rBV	100856	154932	1.17%	0.296%
12	7.248	702	708	711	rBV	120959	197217	1.48%	0.377%
13	7.284	711	714	718	rVV2	174236	230636	1.73%	0.441%
14	7.389	722	732	737	rBV5	131980	283867	2.14%	0.543%
15	7.489	743	749	757	rBV2	923429	1758429	13.23%	3.361%
16	7.777	789	798	803	rBV6	58947	166897	1.26%	0.319%
17	7.842	803	809	816	rBV2	262357	537867	4.05%	1.028%
18	7.924	816	823	828	rBV	364465	628128	4.72%	1.201%
19	7.971	828	831	839	rVB2	136140	247405	1.86%	0.473%
20	8.077	839	849	853	rBV2	213477	450920	3.39%	0.862%
21	8.147	859	861	868	rVB4	94964	144398	1.09%	0.276%
22	8.276	878	883	888	rBV	190999	328962	2.47%	0.629%
23	8.394	898	903	909	rVB2	140784	295068	2.22%	0.564%
24	8.500	916	921	922	rBV2	194720	284869	2.14%	0.545%
25	8.558	927	931	937	rVB2	167448	302036	2.27%	0.577%
26	8.623	937	942	945	rBV2	176327	268178	2.02%	0.513%
27	8.658	945	948	957	rVB	129250	229517	1.73%	0.439%
28	8.811	969	974	977	rBV	99505	157265	1.18%	0.301%
29	8.858	977	982	990	rVB2	129422	250185	1.88%	0.478%
30	8.964	990	1000	1003	rBV3	191761	418147	3.15%	0.799%
31	9.005	1003	1007	1012	rVV	130549	235255	1.77%	0.450%
32	9.087	1012	1021	1030	rVB	912081	1512279	11.38%	2.891%
33	9.181	1030	1037	1038	rBV4	97031	192681	1.45%	0.368%
34	9.205	1038	1041	1047	rVB4	124285	241044	1.81%	0.461%
35	9.281	1047	1054	1058	rBV6	84567	177544	1.34%	0.339%
36	9.645	1108	1116	1121	rVB4	68764	155685	1.17%	0.298%

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
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37	9.734	1121	1131	1136	rBV5	182517	426007	3.20%	0.814%
38	9.792	1136	1141	1144	rBV5	99554	185106	1.39%	0.354%
39	9.933	1156	1165	1170	rVB3	121268	282768	2.13%	0.541%
40	10.004	1170	1177	1181	rBV4	85911	153372	1.15%	0.293%
41	10.074	1186	1189	1193	rVV2	101387	187841	1.41%	0.359%
42	10.186	1201	1208	1213	rBV5	71306	146763	1.10%	0.281%
43	10.268	1218	1222	1229	rVB3	186446	331898	2.50%	0.634%
44	10.638	1275	1285	1291	rBV3	851542	1783366	13.41%	3.409%
45	10.697	1291	1295	1301	rVB4	84185	151555	1.14%	0.290%
46	10.768	1301	1307	1313	rBV4	317722	699626	5.26%	1.337%
47	11.026	1344	1351	1361	rBV	325595	559565	4.21%	1.070%
48	11.150	1364	1372	1379	rVV2	179581	347145	2.61%	0.664%
49	11.672	1456	1461	1466	rBV	203260	343661	2.59%	0.657%
50	11.743	1467	1473	1481	rVB2	226711	418803	3.15%	0.801%
51	11.907	1492	1501	1509	rBV2	201865	387307	2.91%	0.740%
52	12.072	1522	1529	1532	rBV2	317268	540337	4.06%	1.033%
53	12.107	1532	1535	1540	rVB2	176647	241765	1.82%	0.462%
54	12.313	1563	1570	1577	rVB2	377243	660416	4.97%	1.262%
55	12.477	1593	1598	1602	rBV4	130446	224412	1.69%	0.429%
56	12.895	1663	1669	1675	rVB6	70175	145934	1.10%	0.279%
57	13.318	1732	1741	1747	rBV	360552	580623	4.37%	1.110%
58	13.535	1772	1778	1785	rVB5	91827	207502	1.56%	0.397%
59	13.670	1791	1801	1806	rBV6	116857	288788	2.17%	0.552%
60	13.752	1810	1815	1819	rBV2	128163	194762	1.47%	0.372%
61	13.811	1820	1825	1827	rBV2	94670	161762	1.22%	0.309%
62	13.893	1835	1839	1844	rVB	293171	444490	3.34%	0.850%
63	13.976	1844	1853	1857	rBV3	142934	274372	2.06%	0.524%
64	14.123	1873	1878	1882	rBV3	141691	210827	1.59%	0.403%
65	14.181	1884	1888	1892	rVV	229887	300769	2.26%	0.575%
66	14.393	1919	1924	1928	rBV	753197	939274	7.07%	1.795%
67	14.487	1934	1940	1946	rVV2	369031	600749	4.52%	1.148%
68	14.563	1947	1953	1959	rVB6	121737	227753	1.71%	0.435%
69	14.628	1959	1964	1971	rBV	329236	498979	3.75%	0.954%
70	14.692	1971	1975	1985	rVV6	190945	360546	2.71%	0.689%
71	14.886	2002	2008	2013	rVB5	98333	222657	1.67%	0.426%
72	15.021	2026	2031	2034	rBV4	77474	155646	1.17%	0.298%
73	15.116	2042	2047	2055	rVB	651068	1034613	7.78%	1.978%
74	15.251	2065	2070	2078	rVB	335481	545521	4.10%	1.043%
75	15.345	2078	2086	2091	rVB2	354909	583395	4.39%	1.115%
76	15.480	2105	2109	2114	rVB	382005	547691	4.12%	1.047%

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77	15.597	2122	2129	2138	rBV6	202244	389866	2.93%	0.745%
78	15.750	2151	2155	2164	rVB4	149080	264362	1.99%	0.505%
79	15.844	2166	2171	2177	rBV2	138463	227153	1.71%	0.434%
80	16.120	2214	2218	2223	rVB2	168840	228732	1.72%	0.437%
81	16.208	2229	2233	2236	rBV2	366582	524898	3.95%	1.003%
82	16.549	2287	2291	2295	rBV4	96981	148695	1.12%	0.284%
83	16.913	2343	2353	2363	rVV	7589534	13294116	100.00%	25.412%
84	17.002	2364	2368	2372	rVV	328585	454126	3.42%	0.868%
85	17.054	2373	2377	2387	rVB2	189773	332833	2.50%	0.636%
86	17.242	2403	2409	2413	rBV	512533	782395	5.89%	1.496%
87	17.336	2420	2425	2429	rBV	455324	659047	4.96%	1.260%
88	17.530	2453	2458	2464	rVB6	113967	210913	1.59%	0.403%
89	17.736	2489	2493	2499	rVB	273095	359072	2.70%	0.686%
90	17.912	2519	2523	2527	rVB	172259	210412	1.58%	0.402%
91	18.435	2607	2612	2616	rBV	195088	251393	1.89%	0.481%
92	19.064	2716	2719	2721	rBV2	142307	185076	1.39%	0.354%
93	19.634	2810	2816	2822	rBV	553470	883561	6.65%	1.689%
94	20.180	2905	2909	2917	rBV4	103374	179445	1.35%	0.343%
95	21.161	3071	3076	3082	rVB2	236107	331639	2.49%	0.634%
96	21.485	3126	3131	3138	rVB2	505957	792120	5.96%	1.514%
97	22.143	3237	3243	3250	rVB4	81685	152671	1.15%	0.292%
98	22.260	3257	3263	3269	rBV4	94786	185980	1.40%	0.356%
99	24.322	3603	3614	3624	rVB	306610	840201	6.32%	1.606%
100	24.528	3639	3649	3659	rVB	339168	946204	7.12%	1.809%

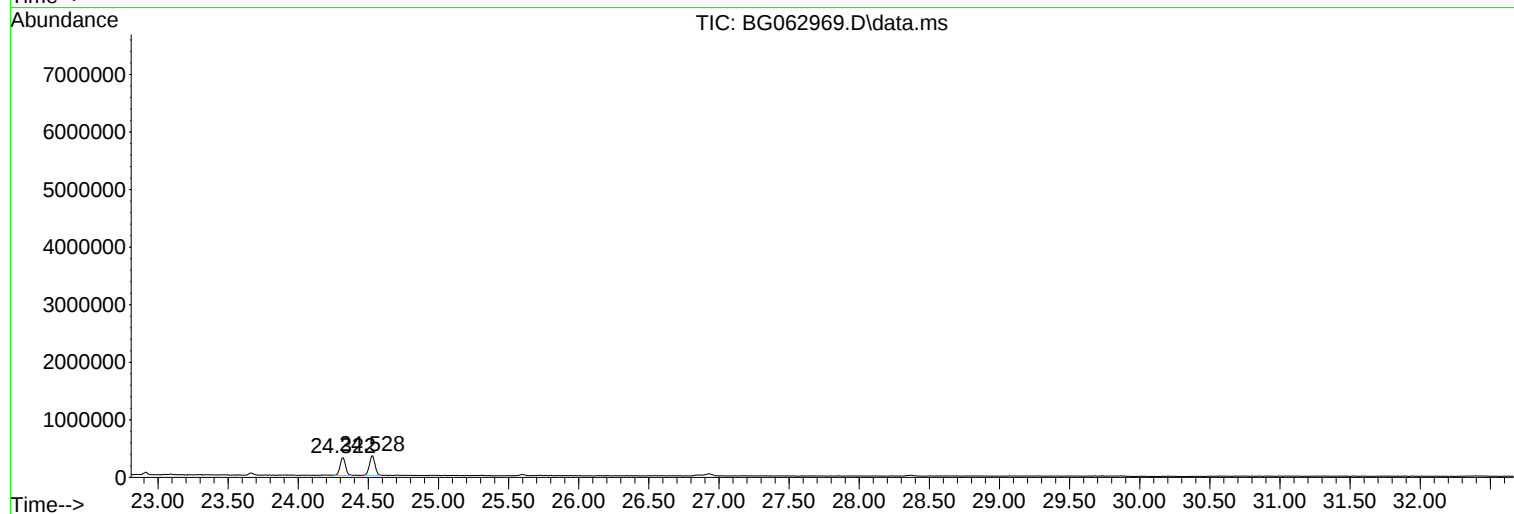
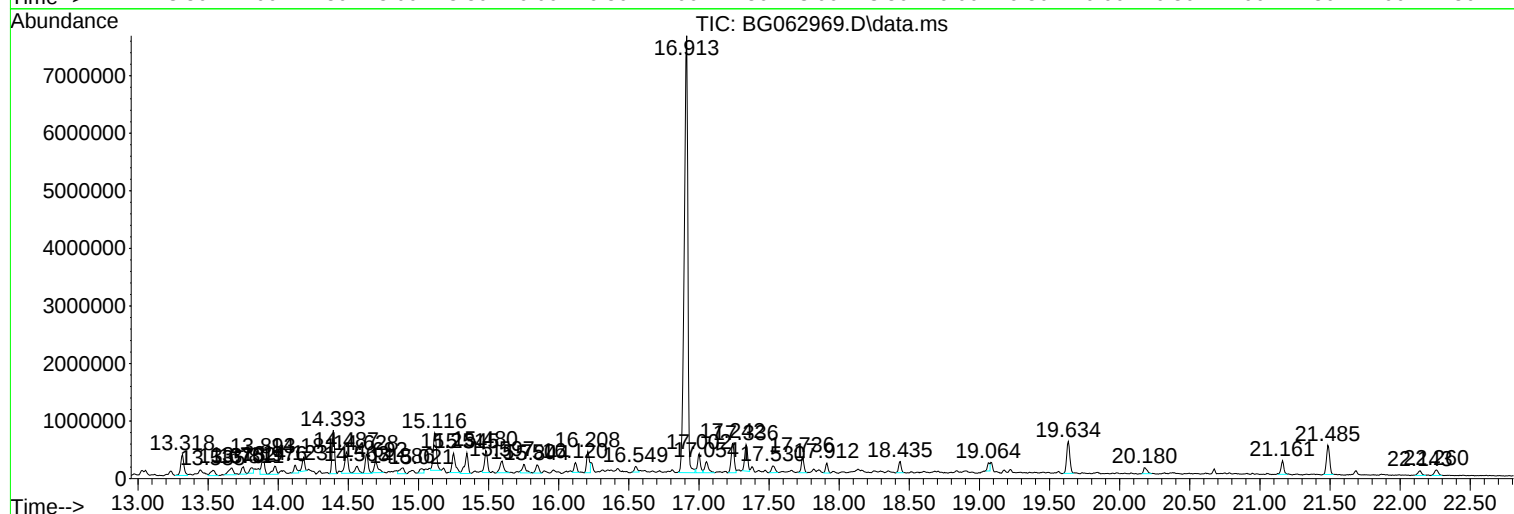
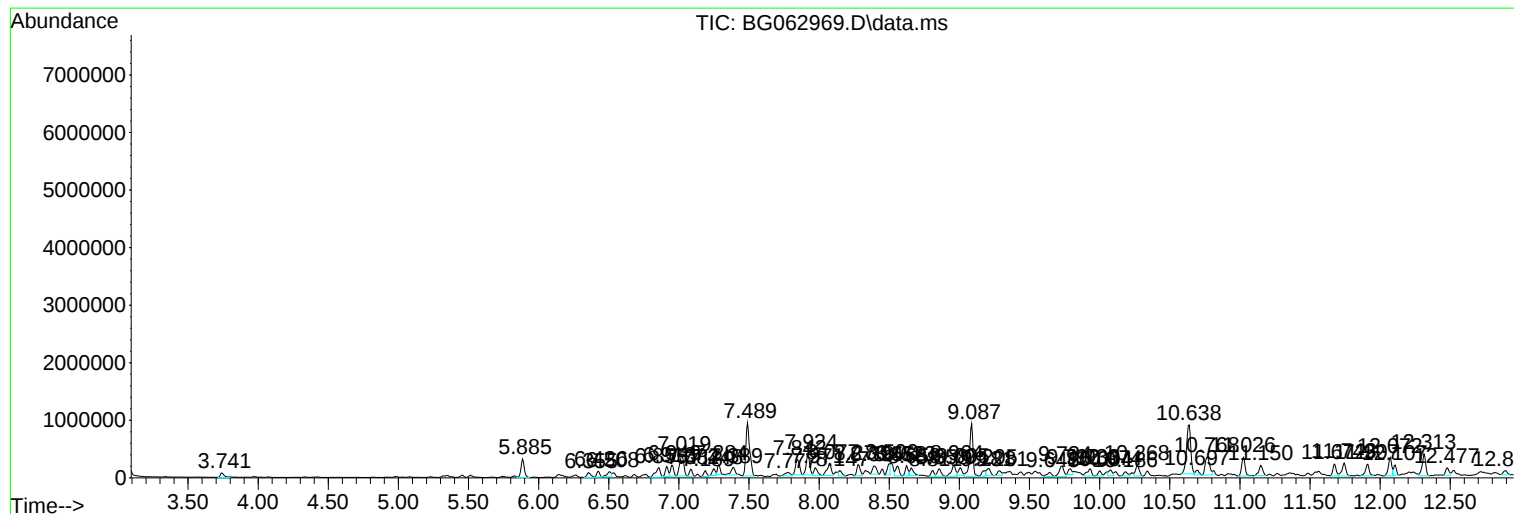
Sum of corrected areas: 52314701

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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



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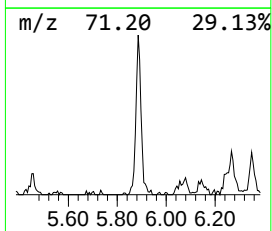
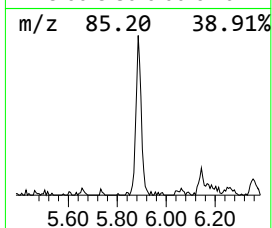
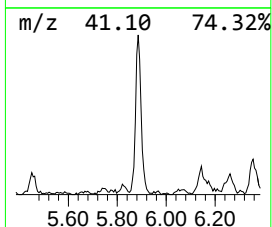
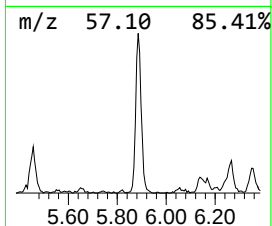
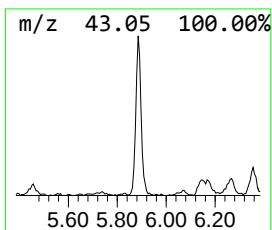
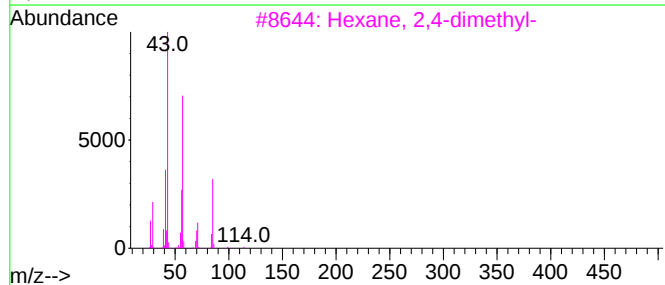
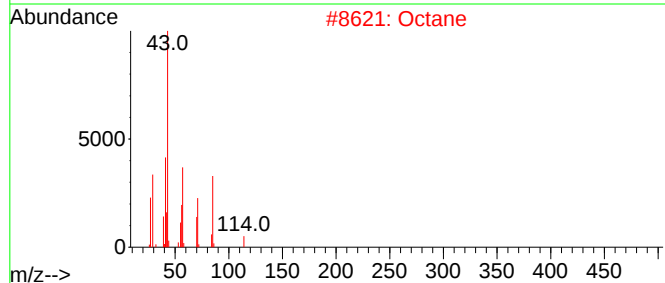
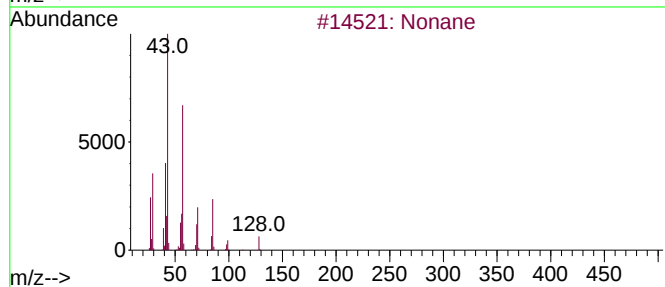
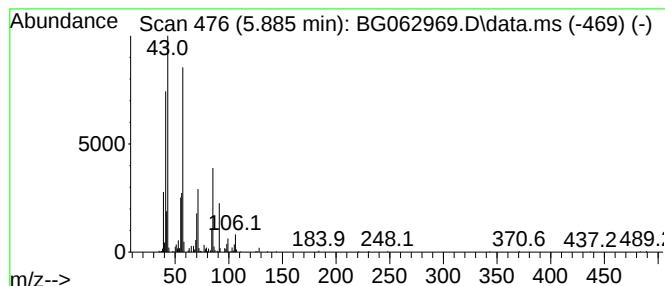
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.885	19.67 ng/ul	528884	1,4-Dichlorobenzene-d4			7.853
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane		128	C9H20	000111-84-2	90
2	Octane		114	C8H18	000111-65-9	64
3	Hexane, 2,4-dimethyl-		114	C8H18	000589-43-5	59
4	4,4-Dimethyl octane		142	C10H22	015869-95-1	59
5	Decane, 1,1'-oxybis-		298	C20H42O	002456-28-2	59



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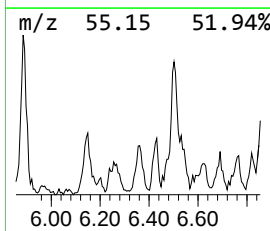
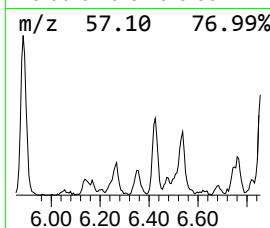
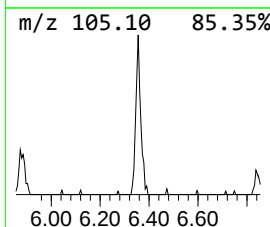
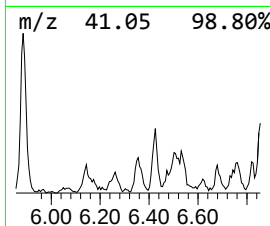
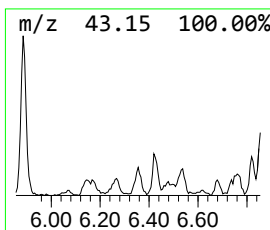
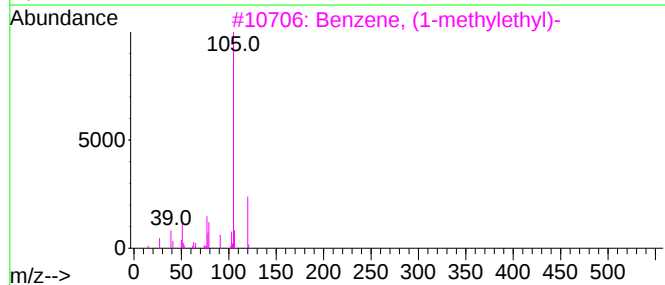
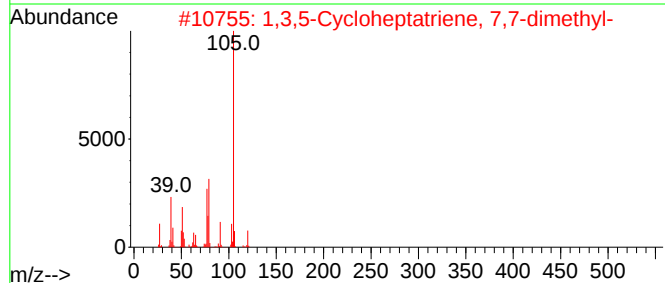
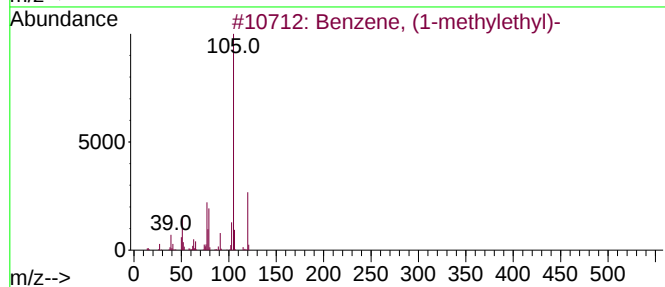
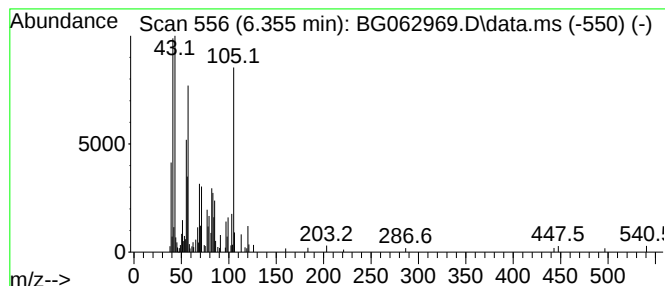
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.355	6.29 ng/ul	169180	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	49
2			1,3,5-Cycloheptatriene, 7,7-dime...	120	C9H12	007557-11-1	43
3			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	43
4			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	43
5			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	43



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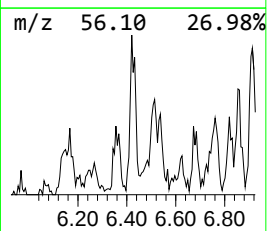
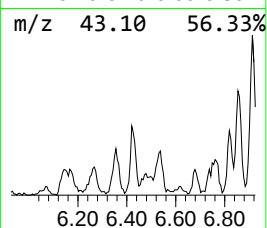
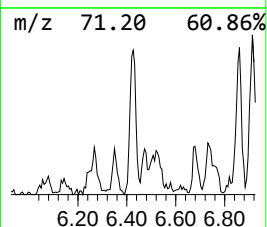
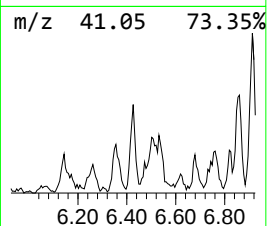
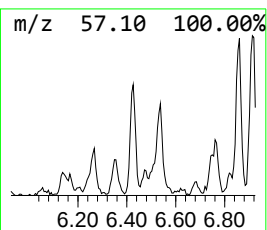
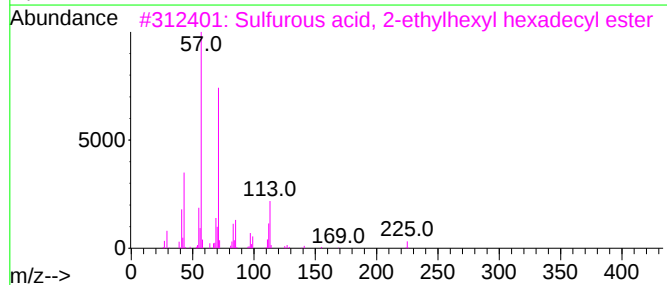
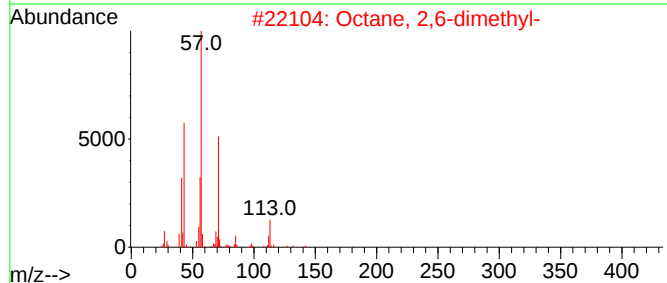
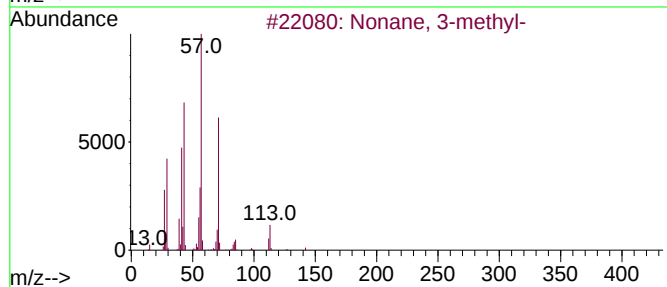
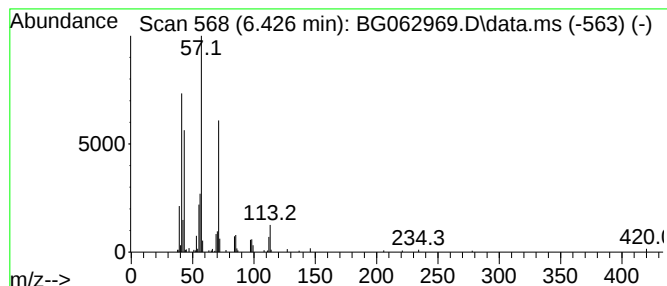
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.426	5.70 ng/ul	153280	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	72
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	72
3			Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	59
4			Sulfurous acid, 2-ethylhexyl pen...	404	C23H48O3S	1000309-19-8	59
5			Sulfurous acid, butyl 2-ethylhex...	250	C12H26O3S	1000309-18-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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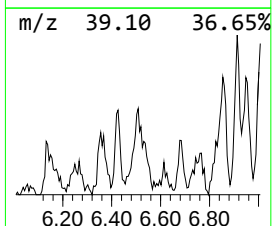
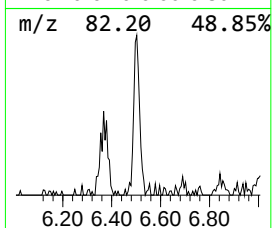
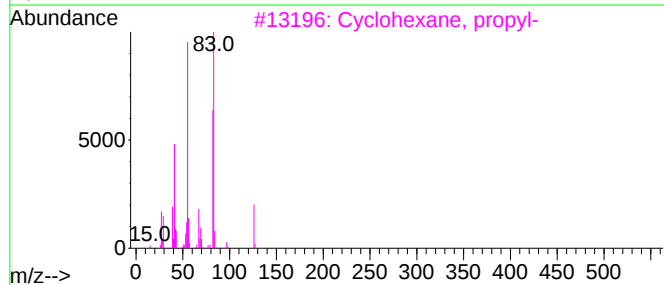
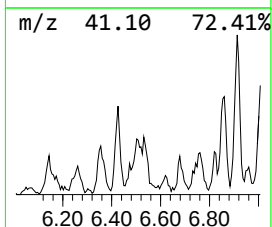
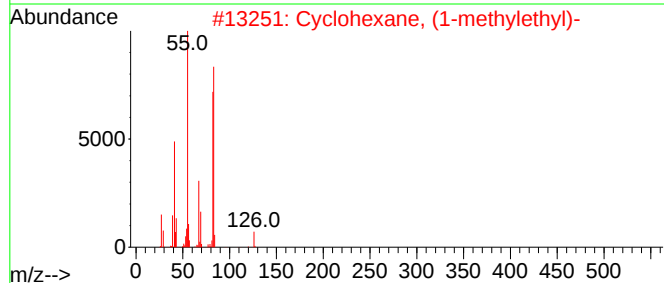
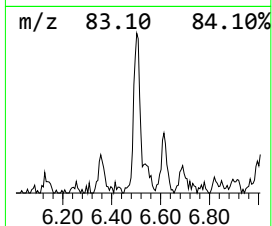
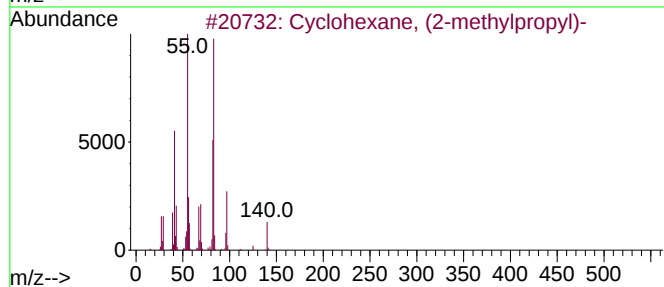
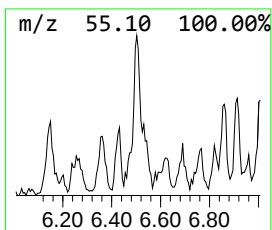
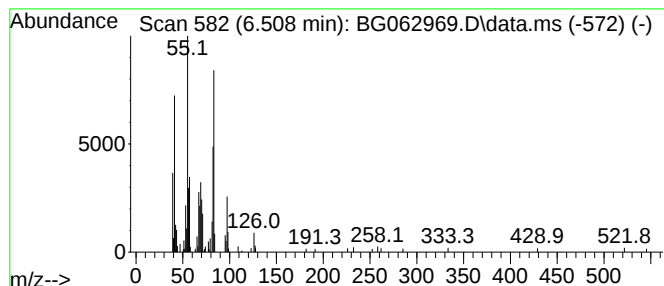
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Cyclic6.508 Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.508	6.84 ng/ul	183853	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	53
2			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	52
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	50
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	47
5			Cyclohexane, methyl-	98	C7H14	000108-87-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
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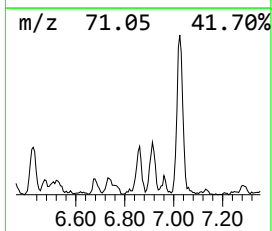
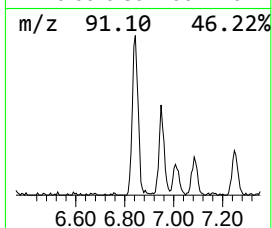
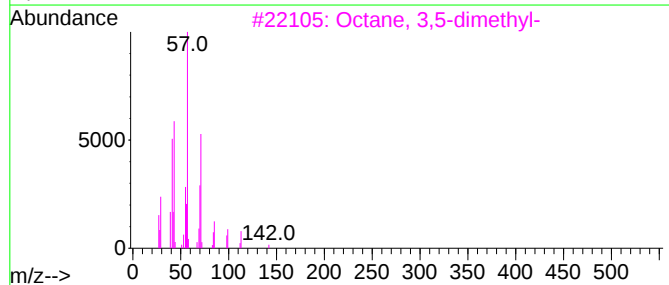
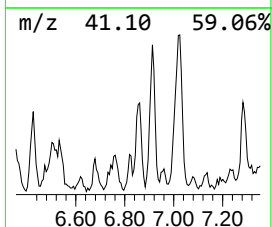
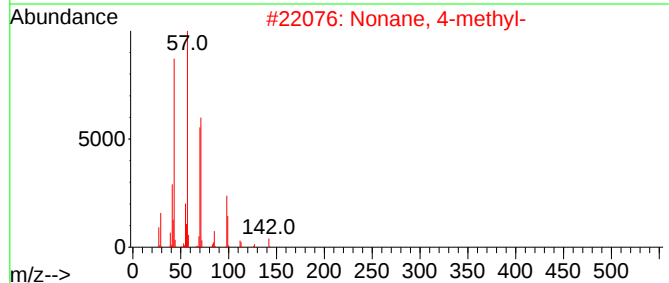
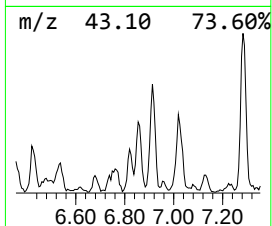
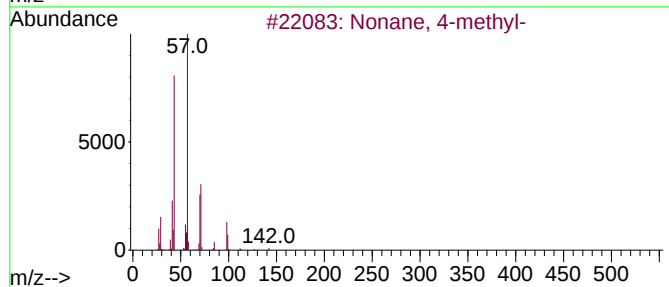
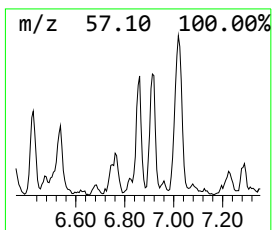
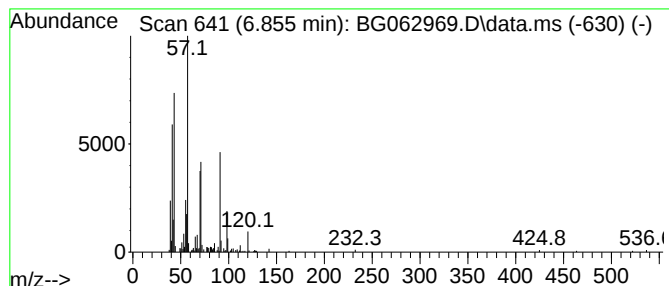
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.855	14.66 ng/ul	394369	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	72
2			Nonane, 4-methyl-	142	C10H22	017301-94-9	64
3			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	64
4			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
5			Undecane, 4,5-dimethyl-	184	C13H28	017312-79-7	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
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ClientSampleId :
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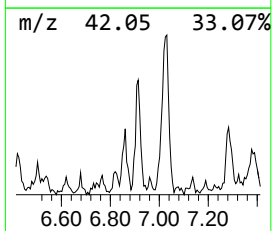
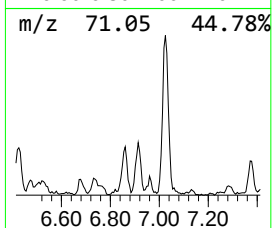
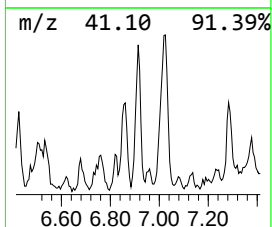
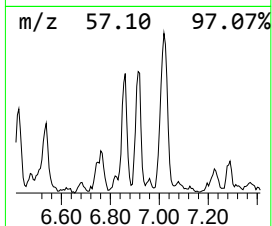
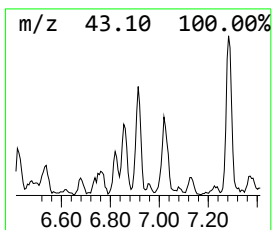
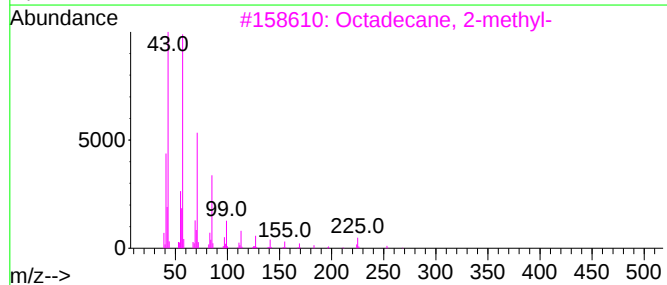
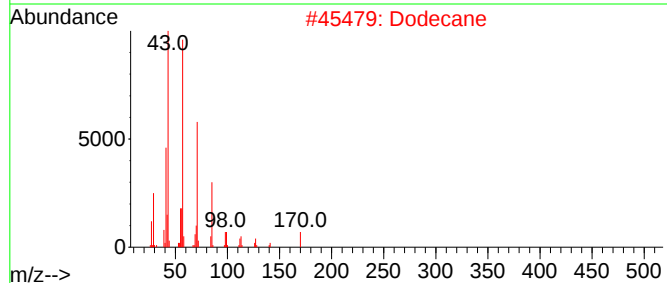
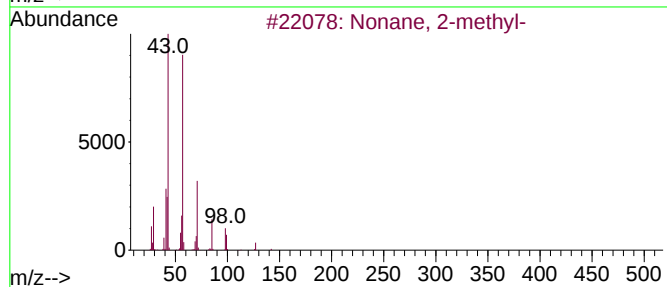
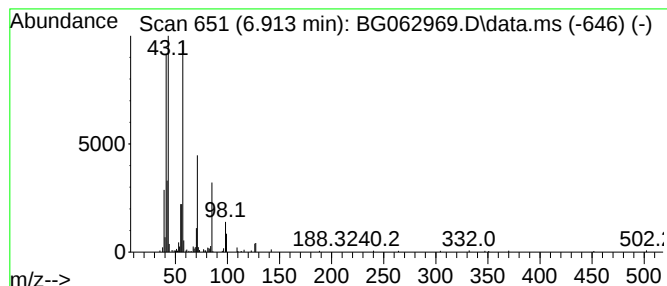
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.913	9.79 ng/ul	263197	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	87
2			Dodecane	170	C12H26	000112-40-3	72
3			Octadecane, 2-methyl-	268	C19H40	001560-88-9	64
4			Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	64
5			Undecane, 5-ethyl-	184	C13H28	017453-94-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
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Sample : P3898-20ME
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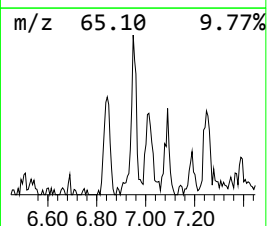
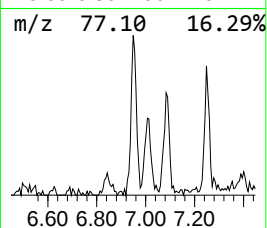
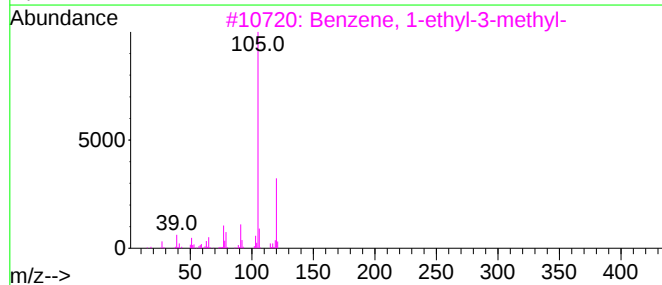
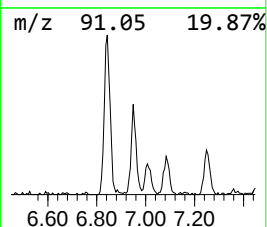
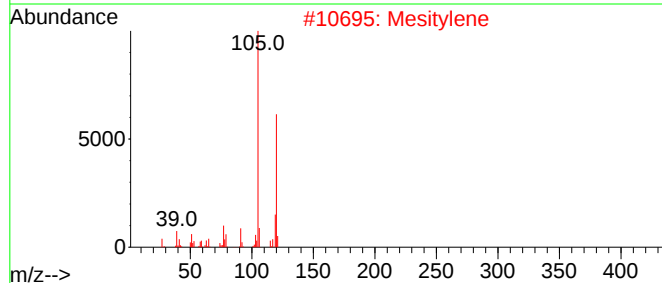
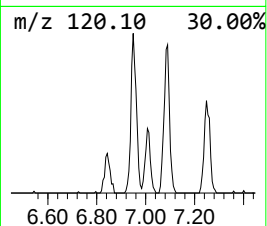
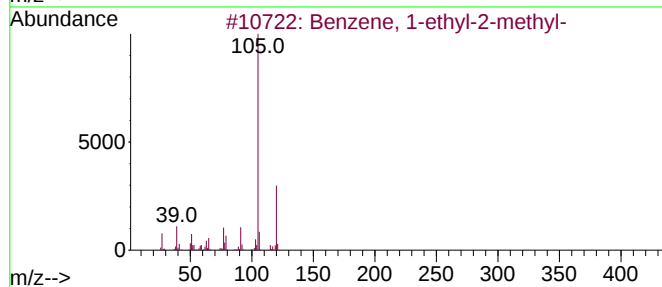
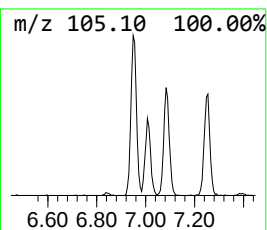
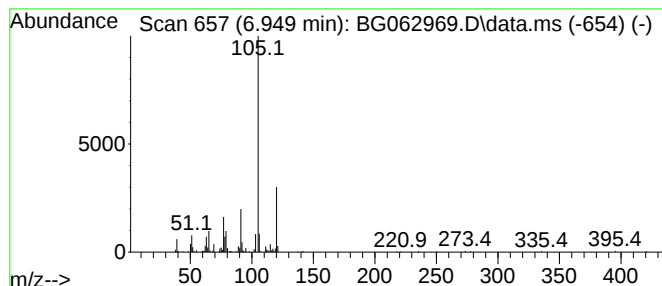
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.949	10.55 ng/ul	283686	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
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Sample : P3898-20ME
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Instrument :
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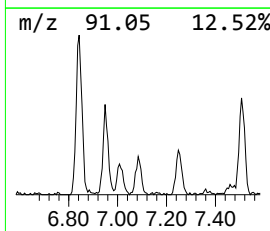
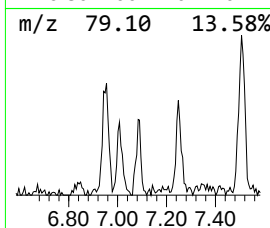
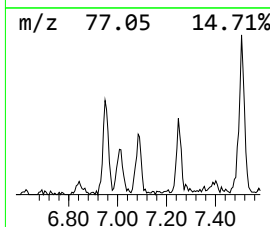
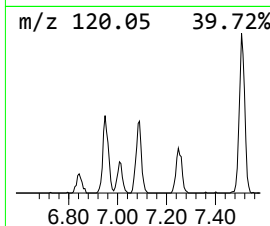
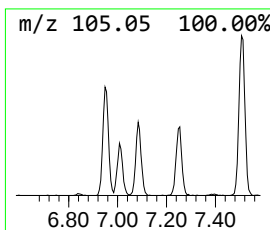
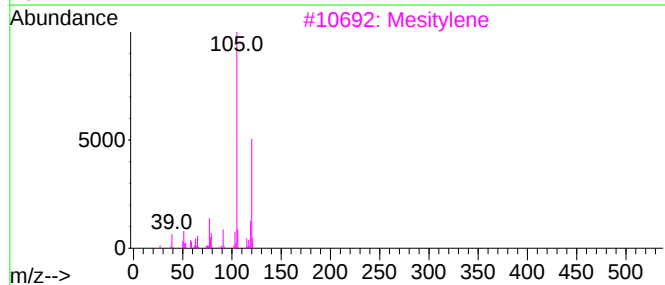
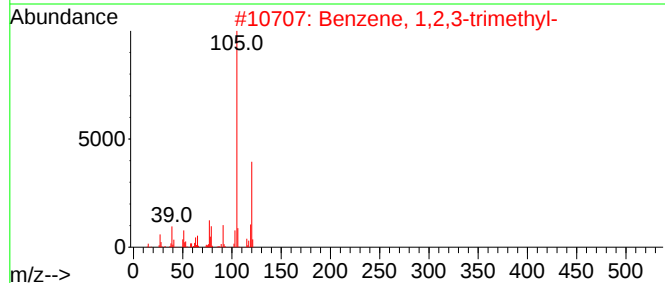
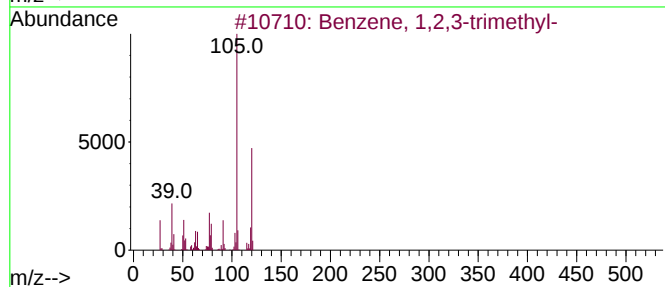
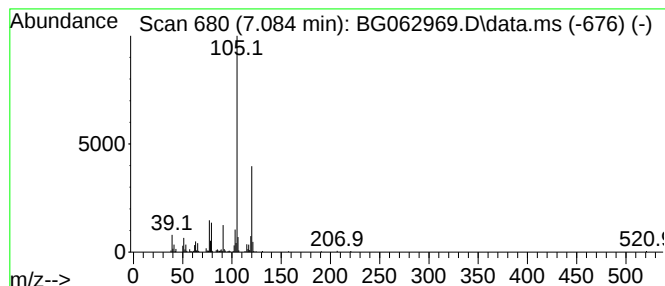
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.084	6.75 ng/ul	181416	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	93
3			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
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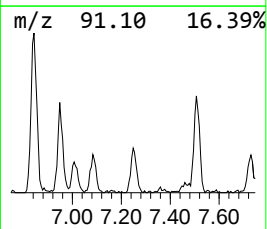
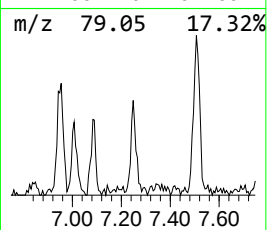
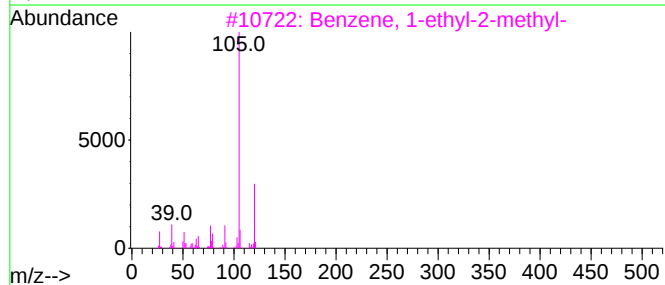
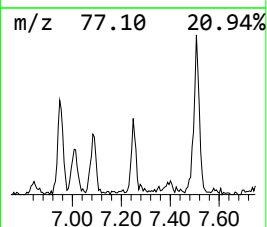
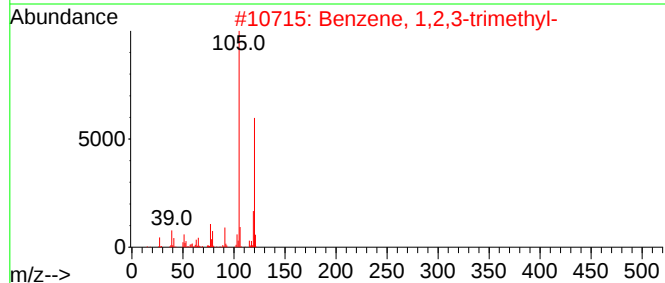
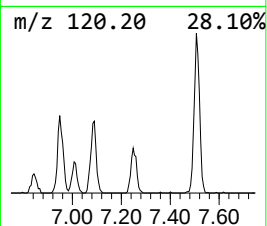
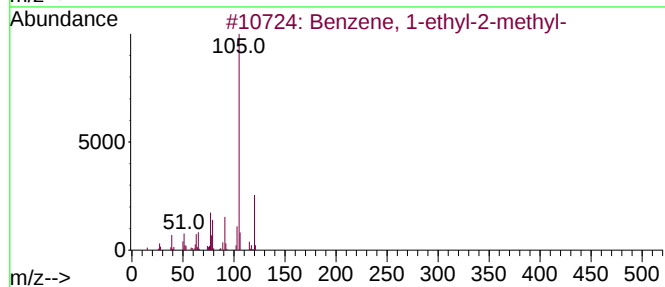
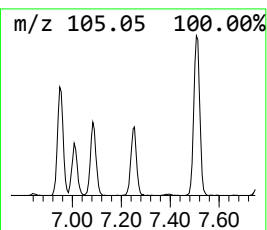
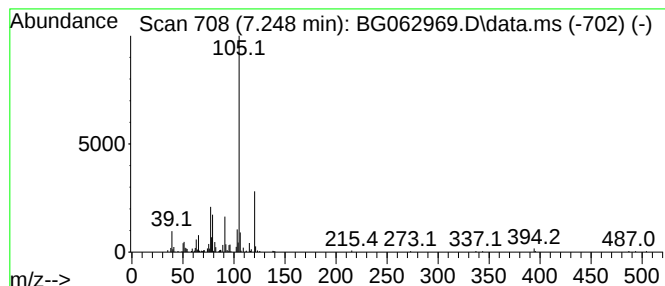
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-ethyl-4-methyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.248	7.33 ng/ul	197217	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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ALS Vial : 33 Sample Multiplier: 1

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ClientSampleId :
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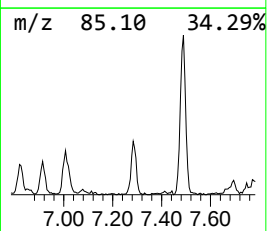
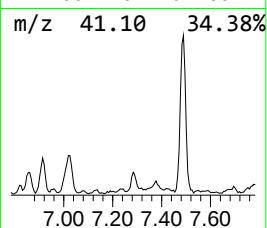
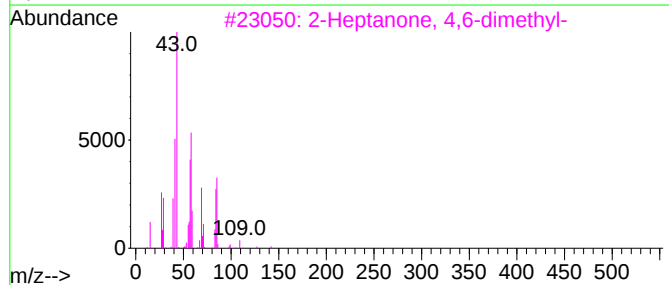
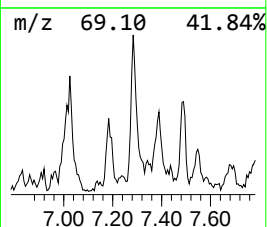
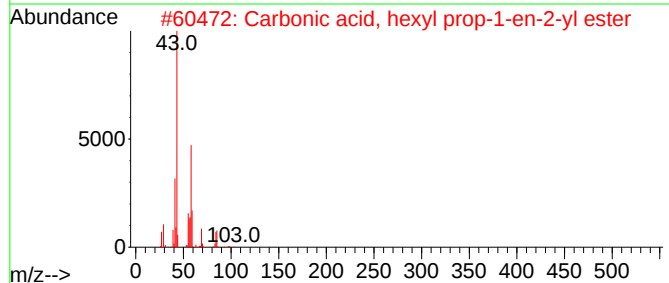
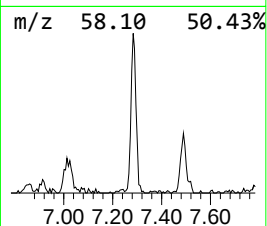
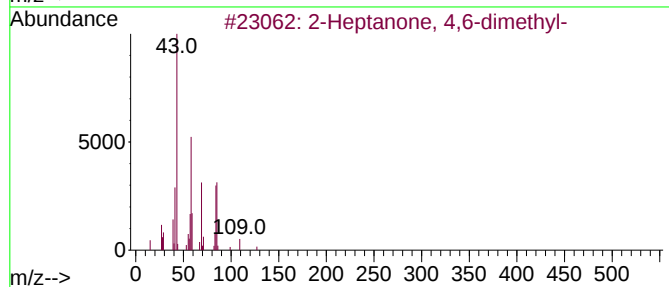
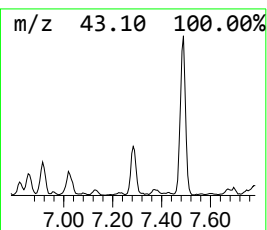
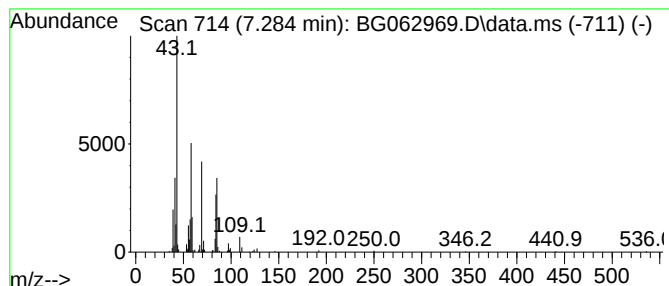
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2-Heptanone, 4,6-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.284	8.58 ng/ul	230636	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	80		
2	Carbonic acid, hexyl prop-1-en-2-yl ester	186	C10H18O3	1000382-54-2	72		
3	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	59		
4	2-Hexanone, 4-methyl-	114	C7H14O	000105-42-0	37		
5	4-Methylpentan-2-yl propyl carbo...	188	C10H20O3	1000372-81-8	32		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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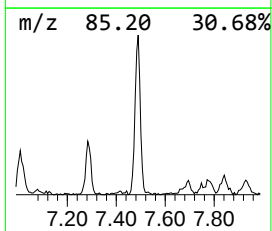
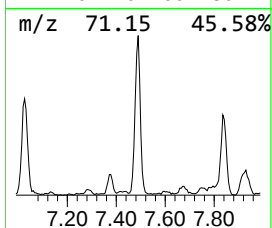
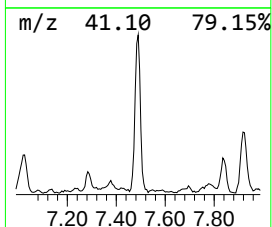
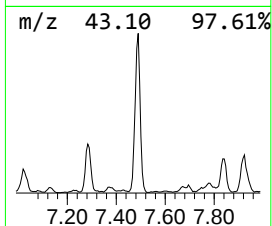
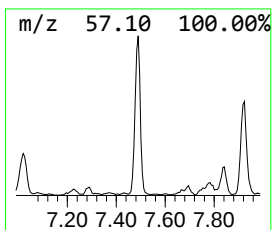
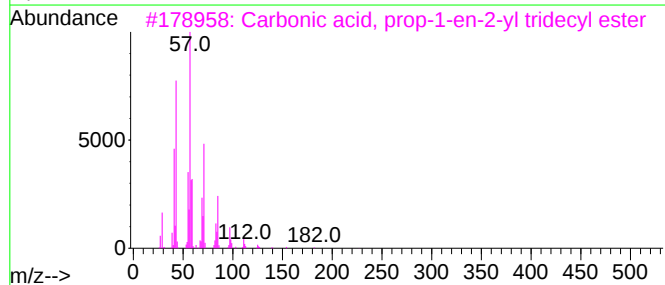
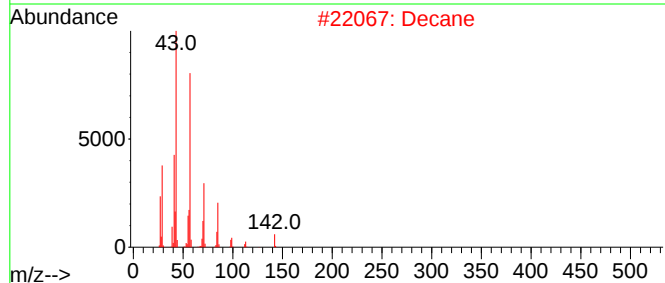
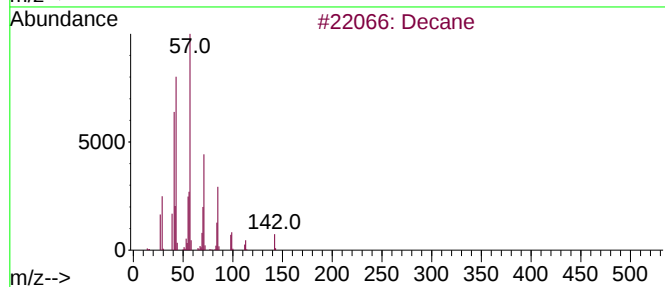
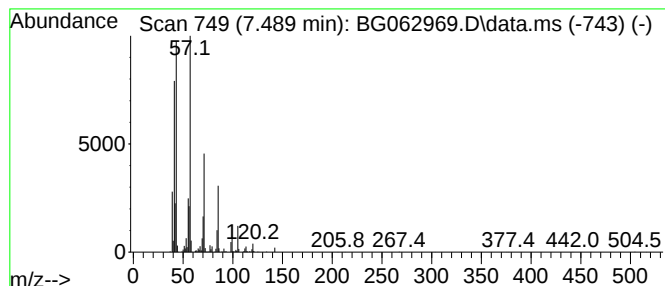
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.489	65.39 ng/ul	1758430	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	90
2			Decane	142	C10H22	000124-18-5	72
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	64
4			Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	64
5			Decane	142	C10H22	000124-18-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

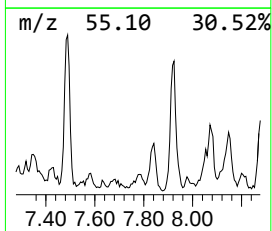
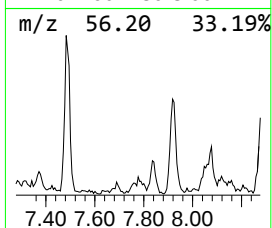
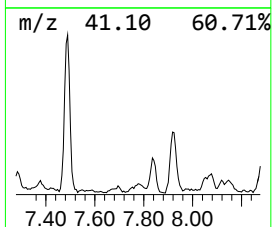
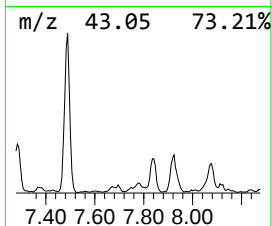
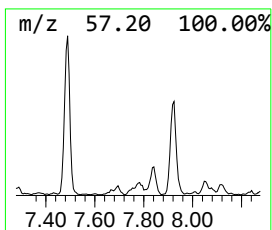
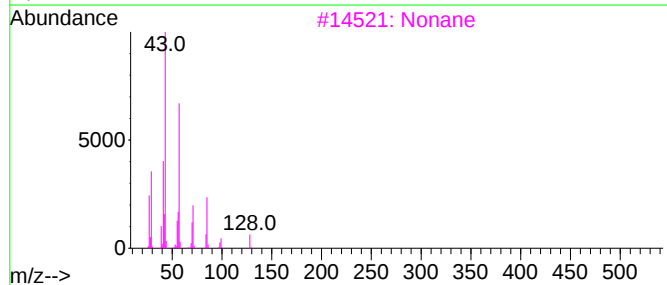
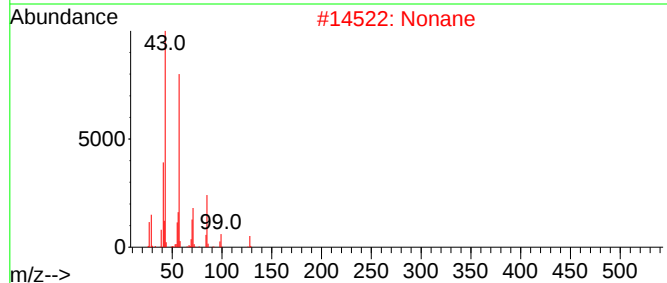
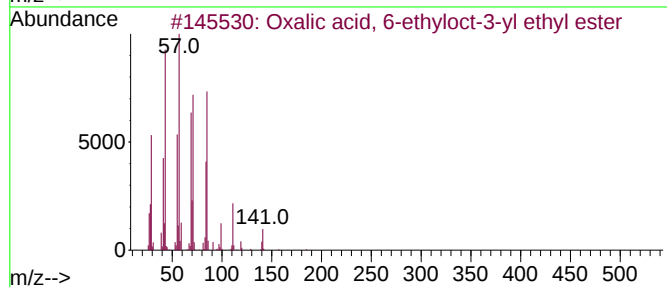
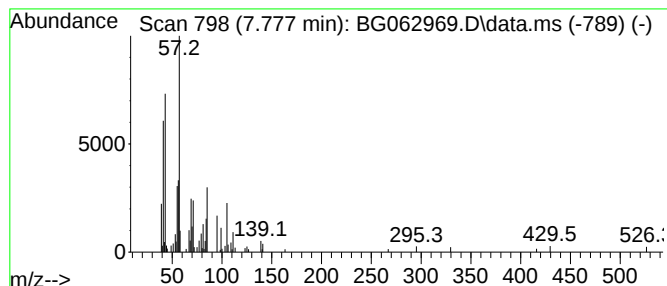
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-02 Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.777	6.21 ng/ul	166897	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 6-ethyloct-3-yl eth...	258	C14H26O4	1000309-33-9	47
2			Nonane	128	C9H20	000111-84-2	43
3			Nonane	128	C9H20	000111-84-2	37
4			Oxalic acid, 6-ethyloct-3-yl hep...	328	C19H36O4	1000309-34-5	35
5			Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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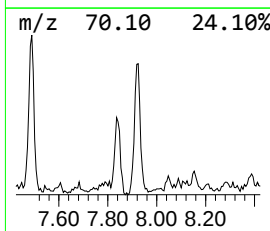
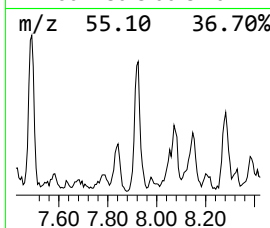
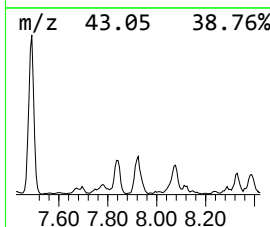
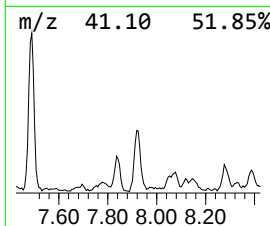
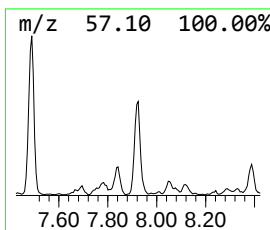
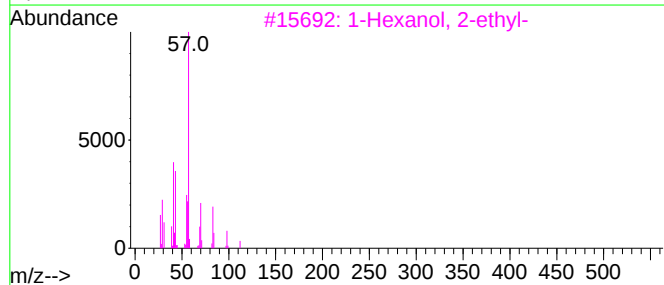
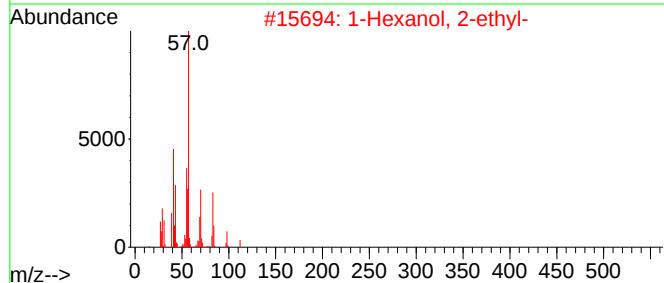
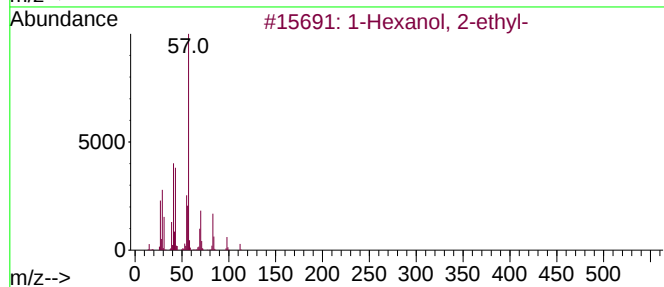
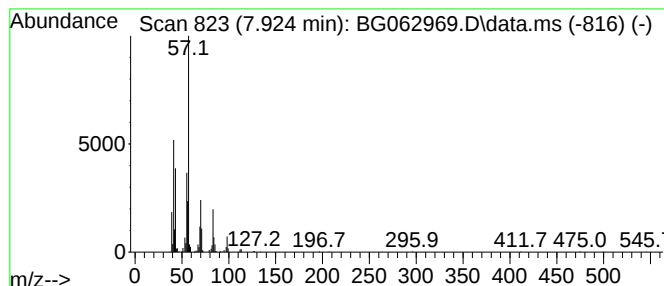
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 1-Hexanol, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.924	23.36 ng/ul	628128	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
5	1-Pentanol, 2-ethyl-4-methyl-			130	C8H18O	000106-67-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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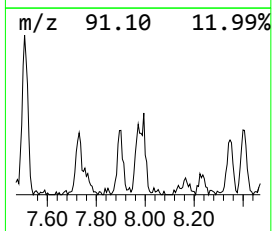
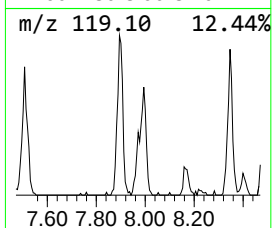
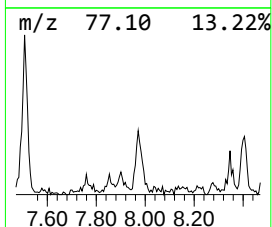
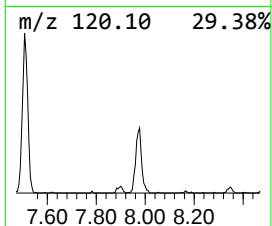
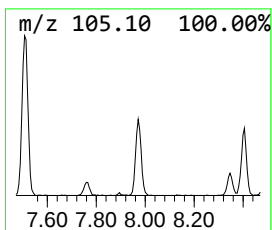
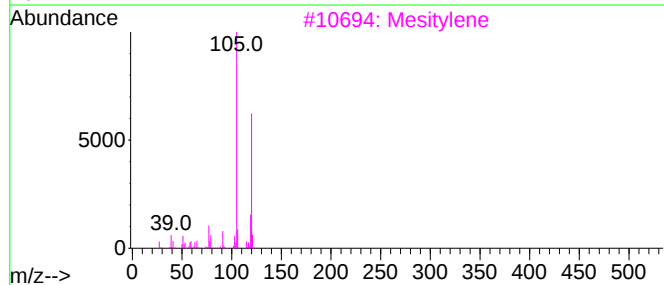
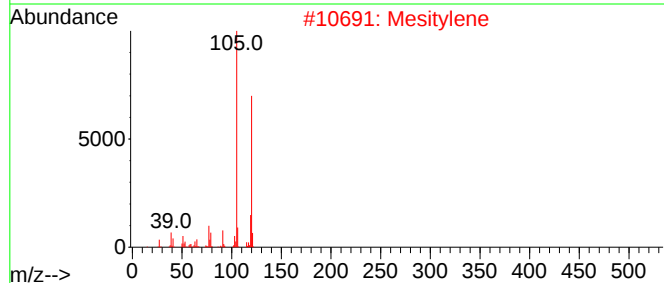
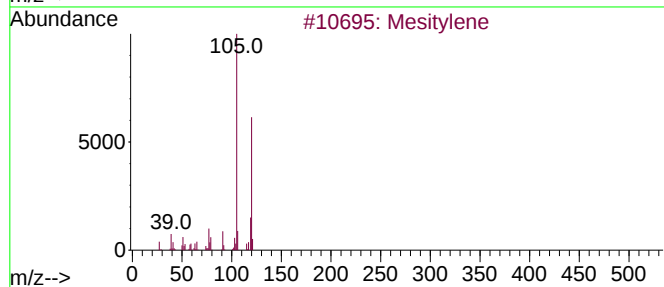
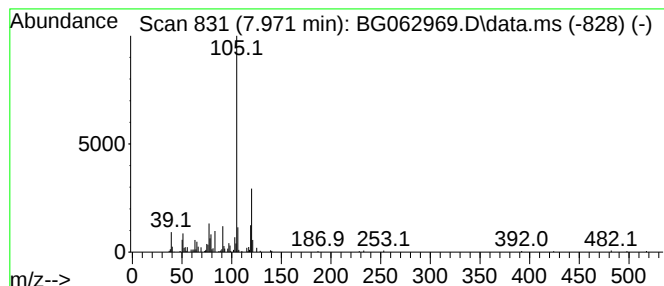
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Mesitylene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.971	9.20 ng/ul	247405	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	81
2			Mesitylene	120	C9H12	000108-67-8	81
3			Mesitylene	120	C9H12	000108-67-8	81
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	81
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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Sample : P3898-20ME
Misc :
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Instrument :
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ClientSampleId :
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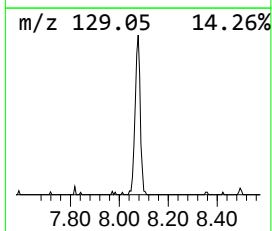
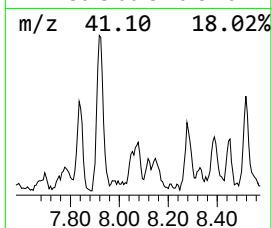
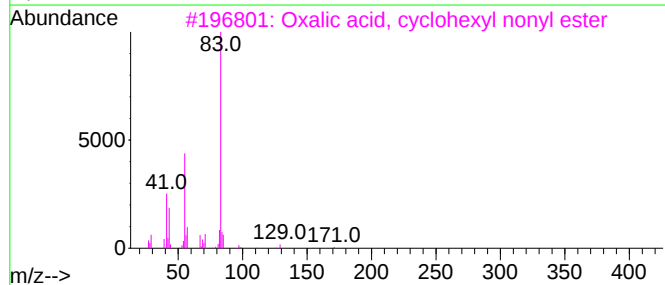
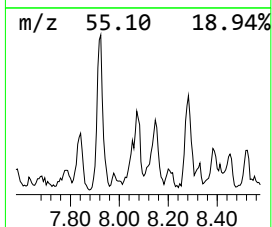
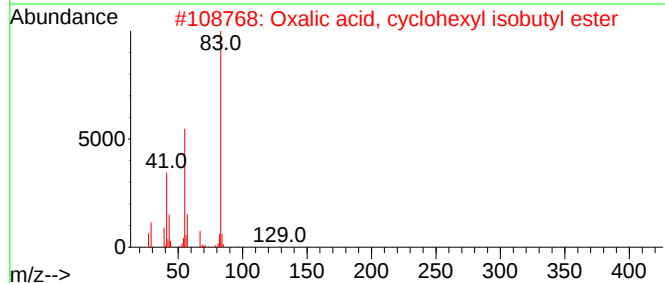
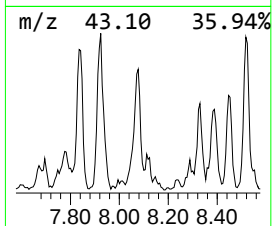
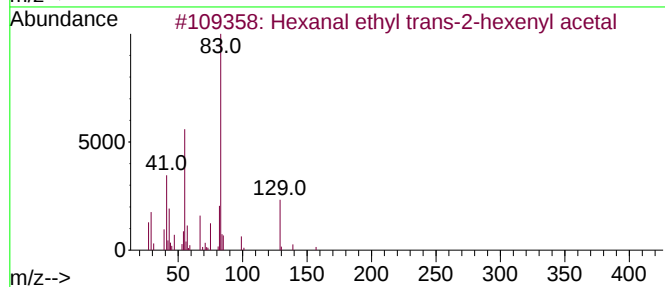
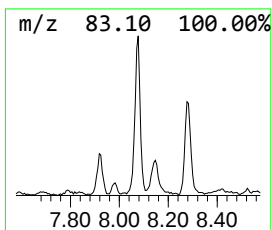
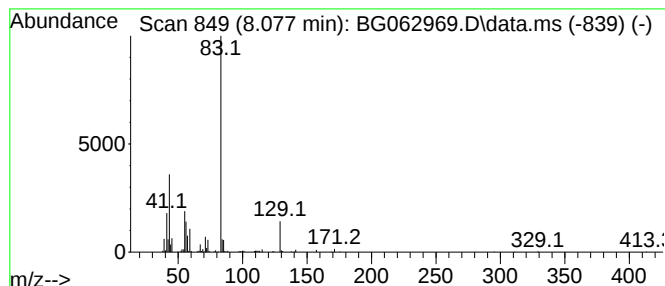
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Hexanal ethyl trans-2-hexen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.077	16.77 ng/ul	450920	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanal ethyl trans-2-hexenyl ac...	228	C14H28O2	1000431-42-6	59
2			Oxalic acid, cyclohexyl isobutyl...	228	C12H20O4	1000309-30-4	42
3			Oxalic acid, cyclohexyl nonyl ester	298	C17H30O4	1000309-31-1	42
4			Oxalic acid, cyclohexyl octyl ester	284	C16H28O4	1000309-31-0	42
5			Oxalic acid, cyclohexyl pentyl e...	242	C13H22O4	1000309-30-6	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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Sample : P3898-20ME
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ALS Vial : 33 Sample Multiplier: 1

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ClientSampleId :
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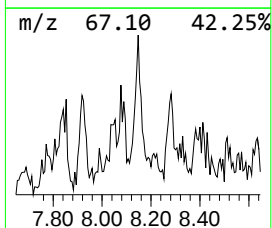
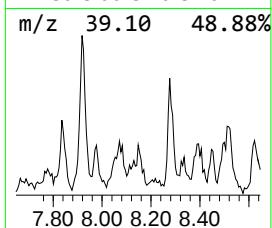
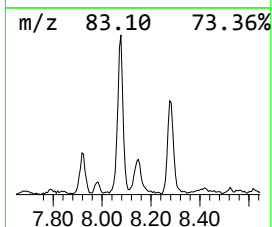
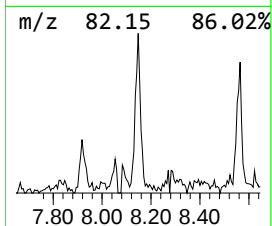
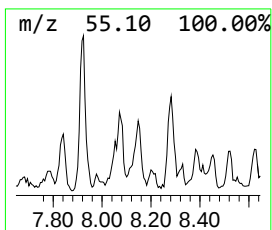
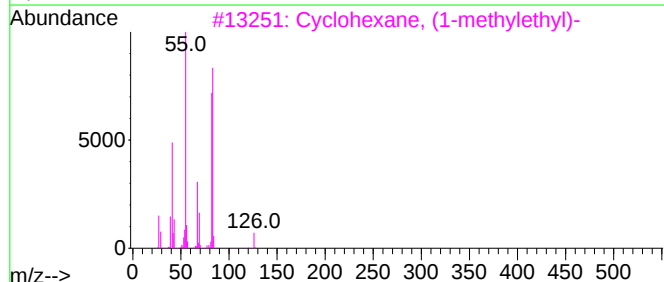
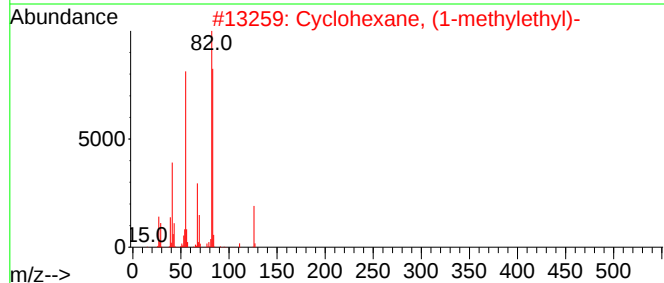
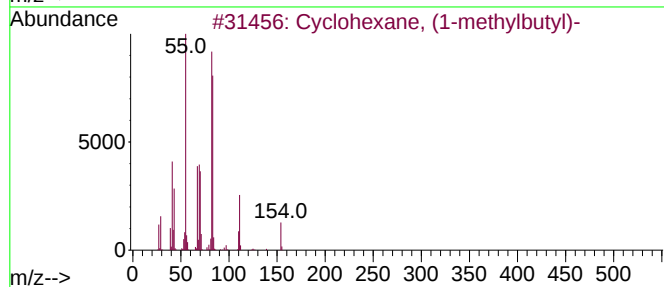
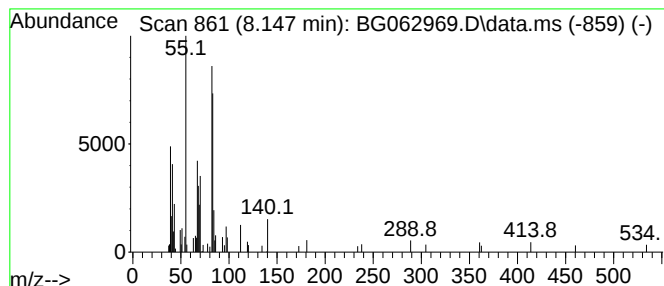
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Cyclic8.147 Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.147	5.37 ng/ul	144398	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylbutyl)-	154	C11H22	061208-94-4	64
2			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	64
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	59
4			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	50
5			Decane, 2-cyclohexyl-	224	C16H32	013151-73-0	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
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Instrument :
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ClientSampleId :
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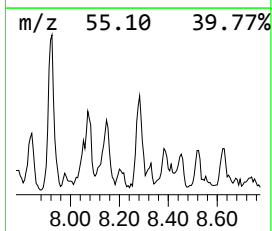
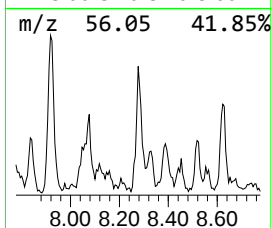
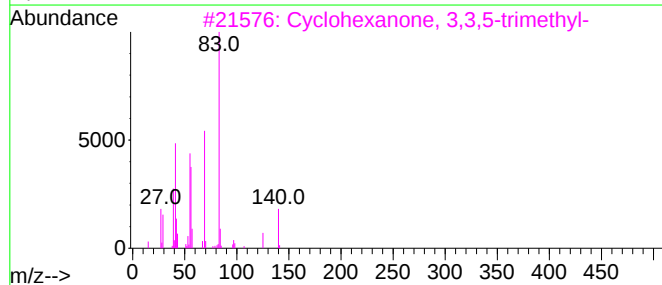
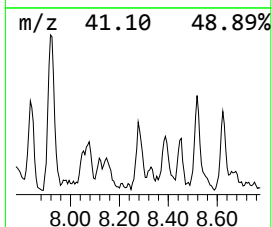
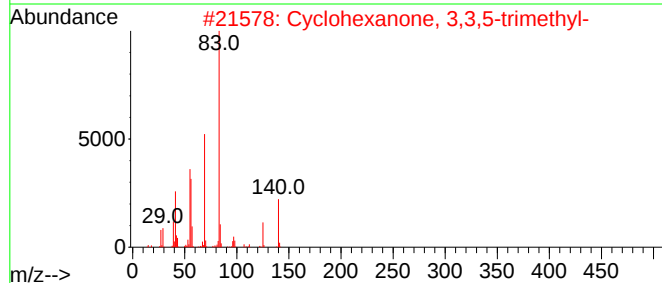
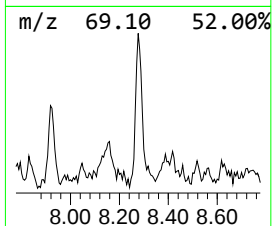
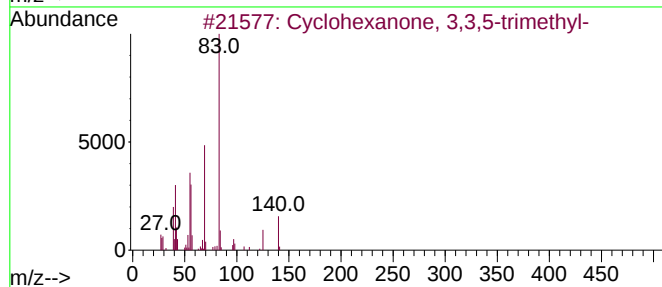
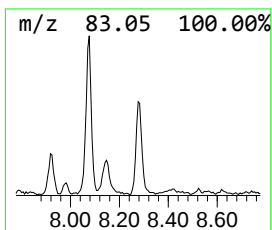
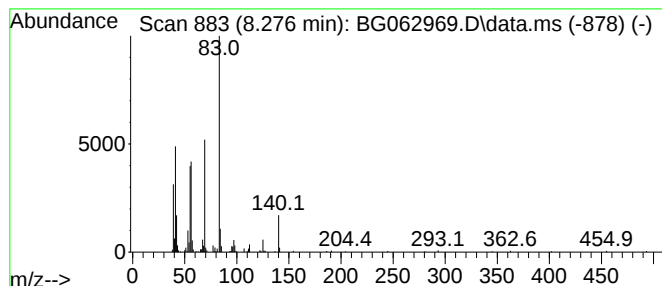
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Cyclohexanone, 3,3,5-trimet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.276	12.23 ng/ul	328962	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
5			1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
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Sample : P3898-20ME
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ClientSampleId :
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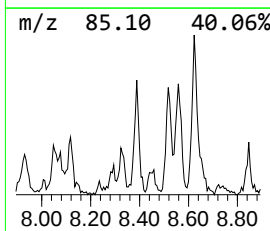
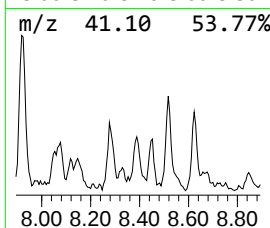
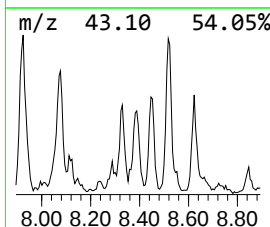
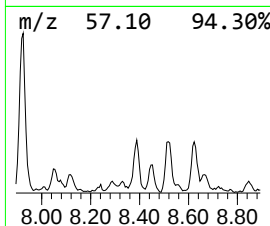
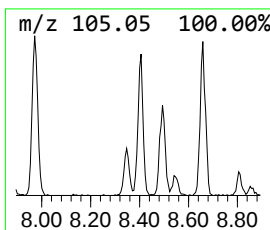
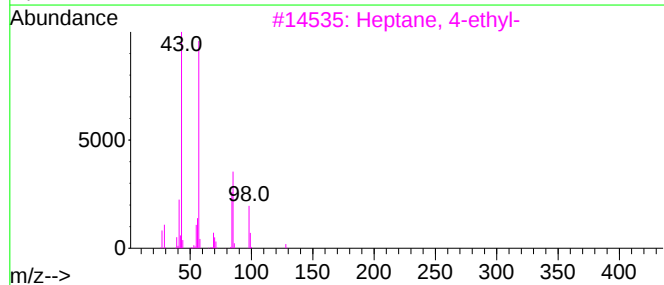
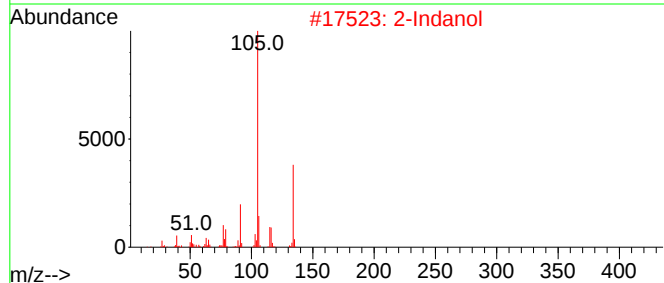
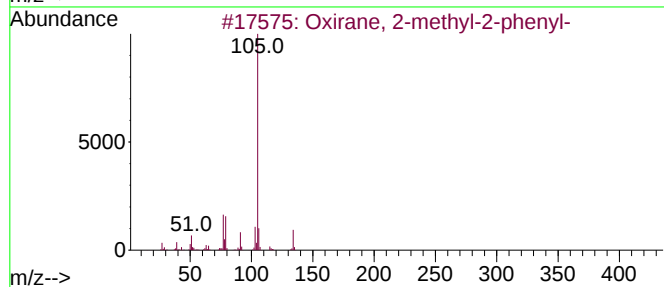
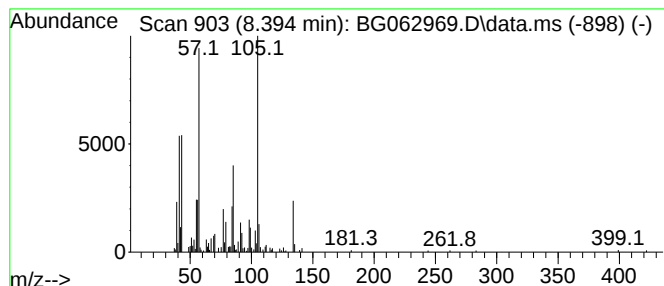
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-03 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.394	10.97 ng/ul	295068	1,4-Dichlorobenzene-d4	7.853

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	46
2		2-Indanol	134	C9H10O	004254-29-9	38
3		Heptane, 4-ethyl-	128	C9H20	002216-32-2	38
4		Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	38
5		Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
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Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

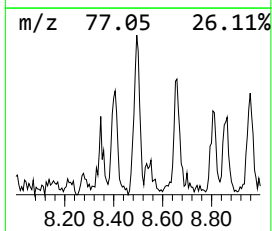
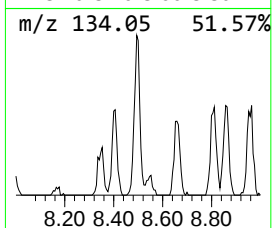
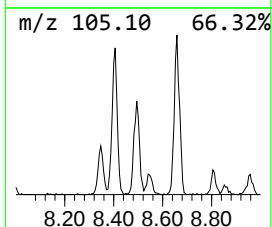
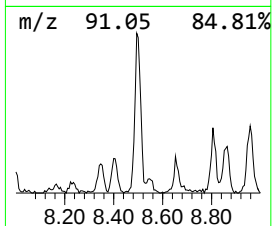
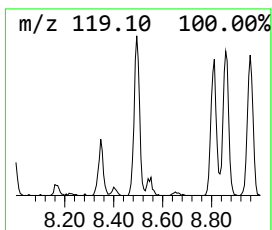
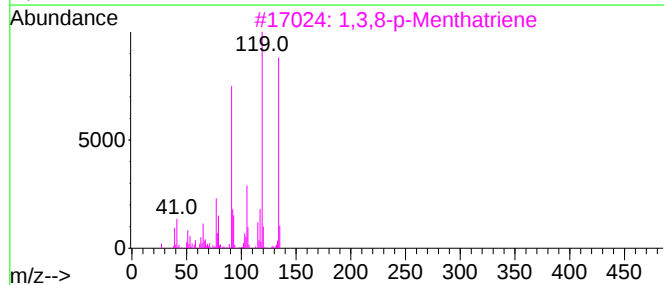
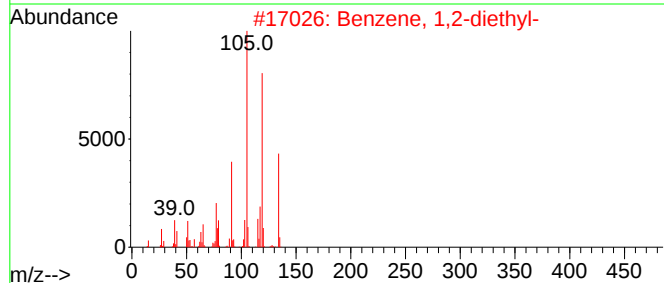
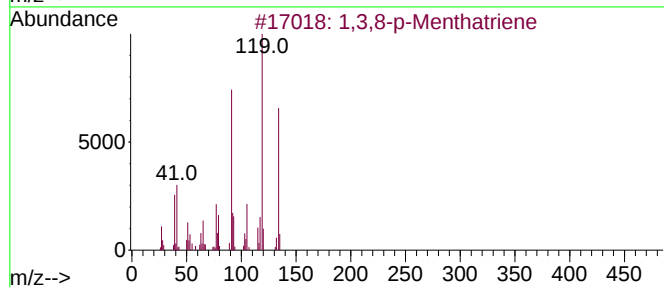
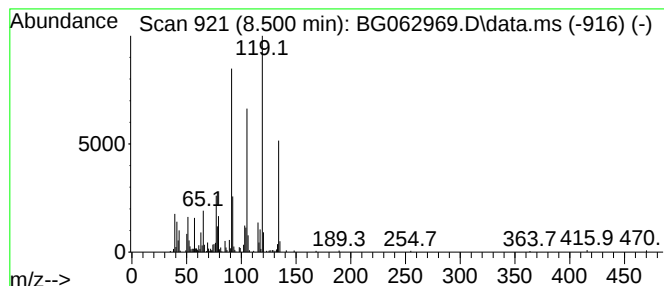
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 1,3,8-p-Menthatriene Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.500	10.59 ng/ul	284869	1,4-Dichlorobenzene-d4			7.853
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1,3,8-p-Menthatriene		134	C10H14	018368-95-1	91
2	Benzene, 1,2-diethyl-		134	C10H14	000135-01-3	89
3	1,3,8-p-Menthatriene		134	C10H14	018368-95-1	87
4	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	81
5	Benzene, 2-ethyl-1,4-dimethyl-		134	C10H14	001758-88-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
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Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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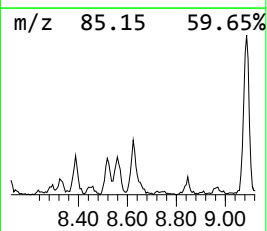
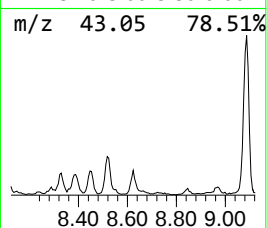
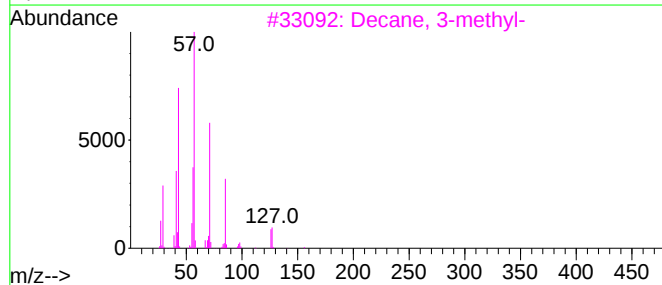
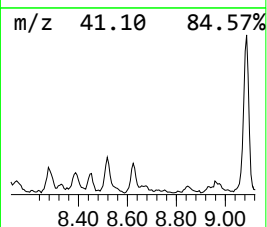
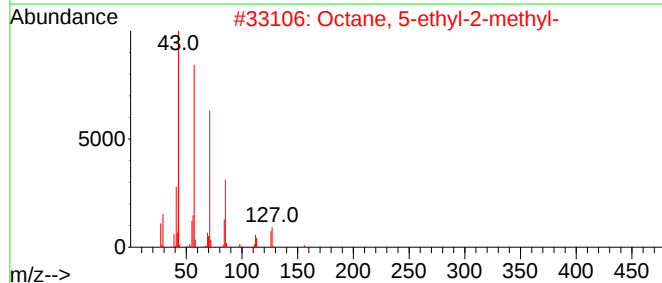
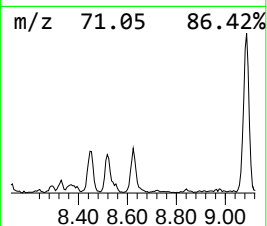
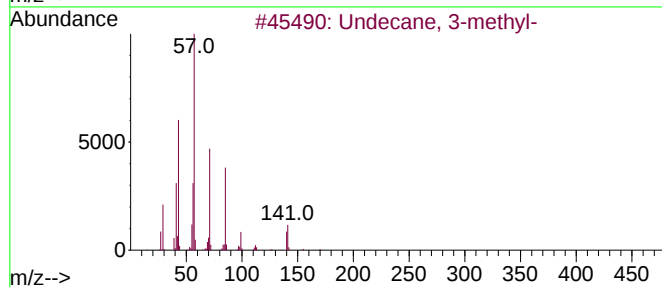
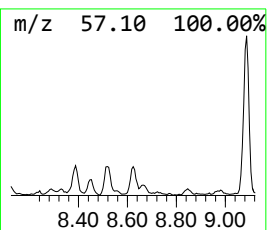
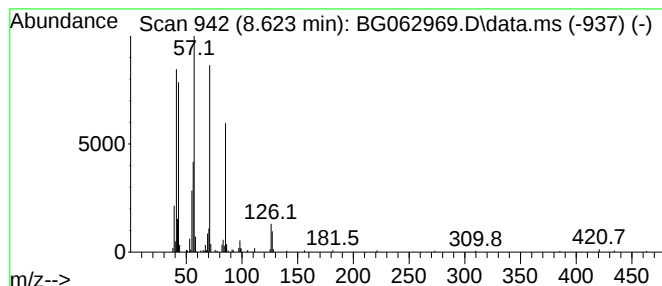
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.623	9.97 ng/ul	268178	1,4-Dichlorobenzene-d4	7.853

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 3-methyl-	170	C12H26	001002-43-3	72
2		Octane, 5-ethyl-2-methyl-	156	C11H24	062016-18-6	64
3		Decane, 3-methyl-	156	C11H24	013151-34-3	64
4		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	59
5		Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
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Sample : P3898-20ME
Misc :
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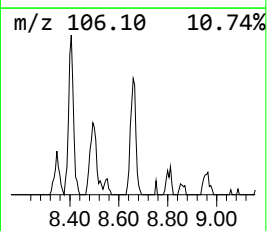
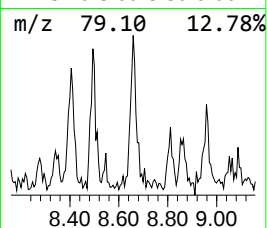
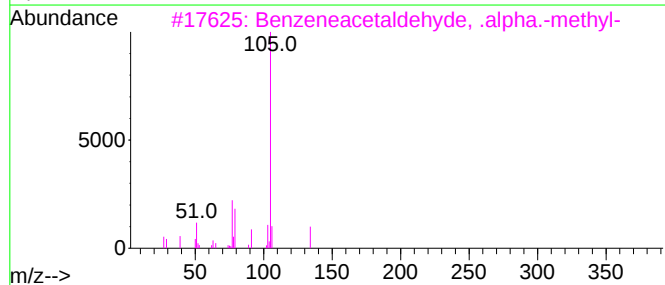
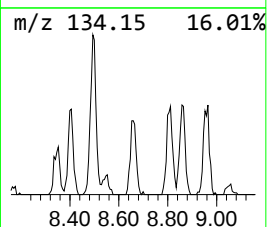
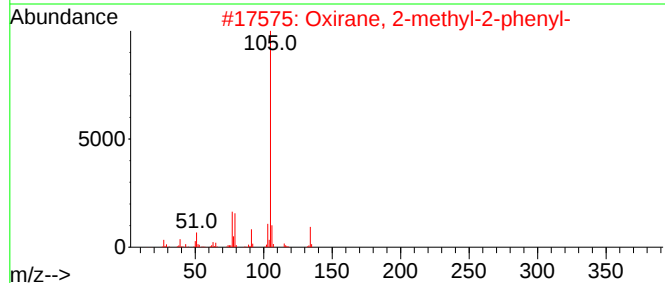
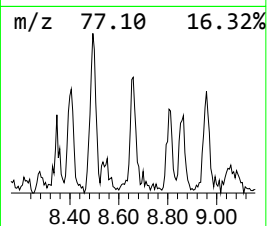
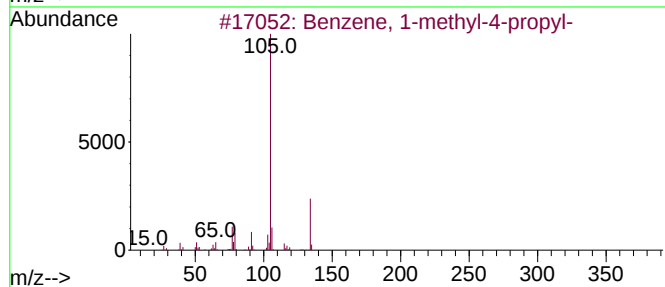
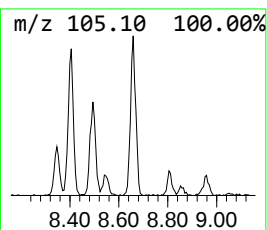
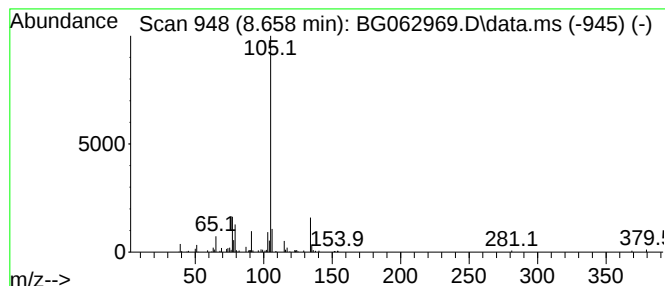
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Benzene, 1-methyl-4-propyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.658	8.53 ng/ul	229517	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
2			Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	91
3			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	87
4			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	86
5			Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
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Acq On : 20 Sep 2024 8:39
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Misc :
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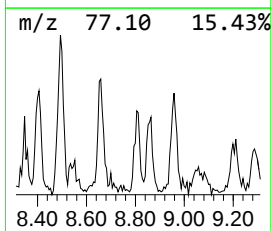
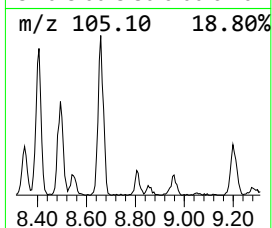
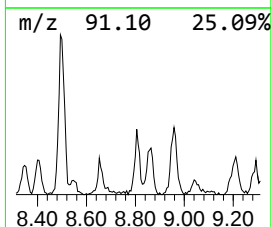
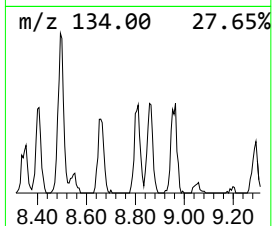
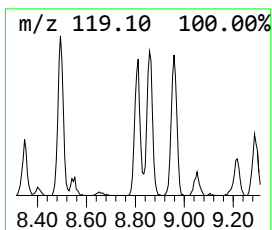
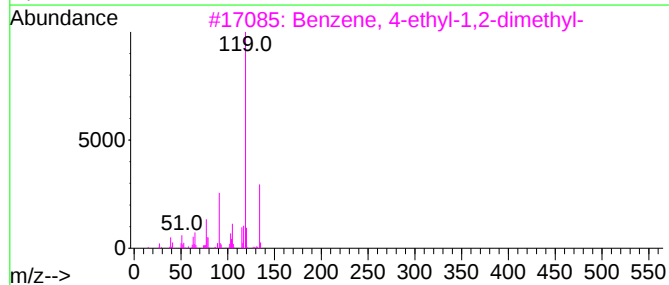
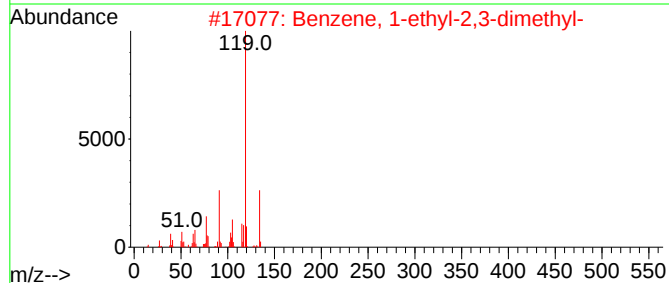
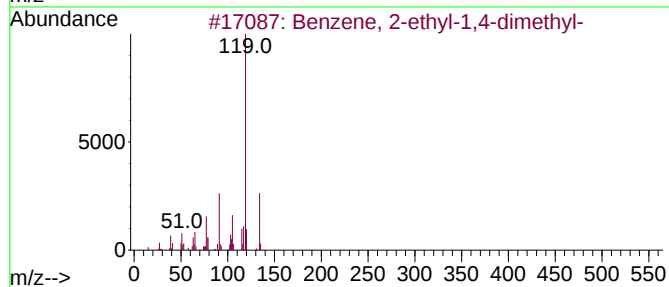
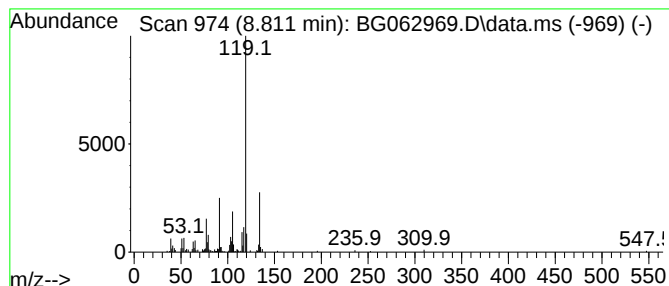
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.811	5.85 ng/ul	157265	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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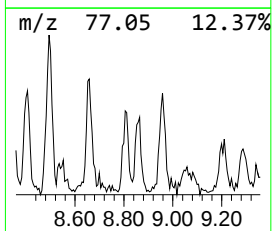
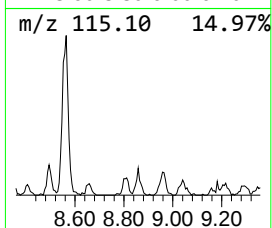
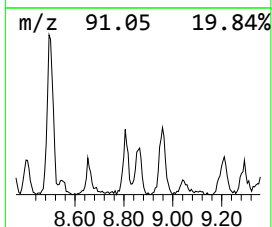
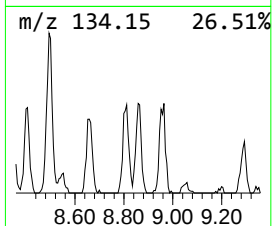
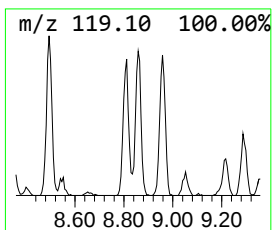
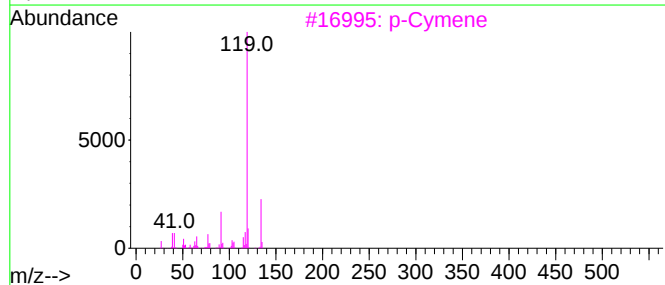
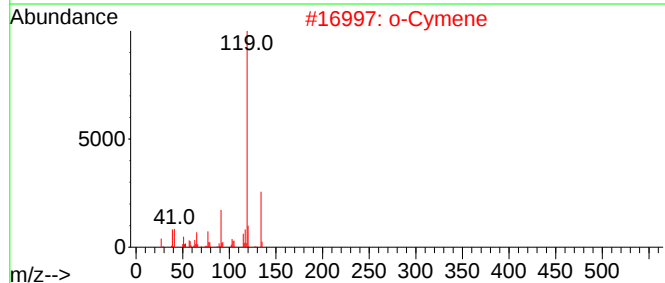
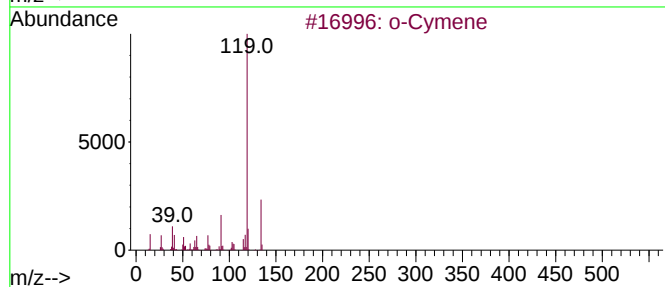
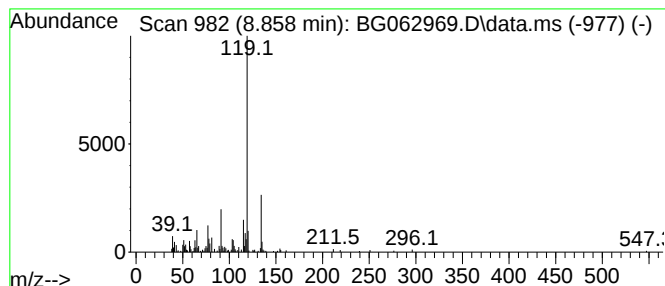
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 o-Cymene Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.858	9.30 ng/ul	250185	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	93
2			o-Cymene	134	C10H14	000527-84-4	93
3			p-Cymene	134	C10H14	000099-87-6	93
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	90
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

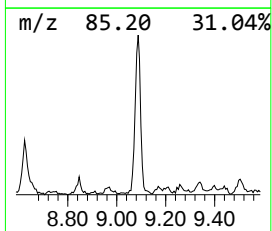
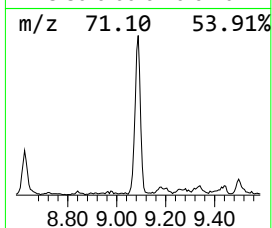
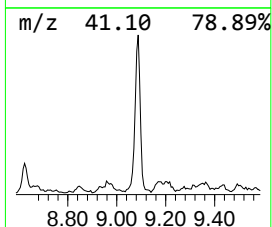
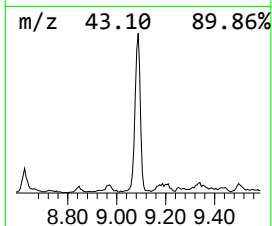
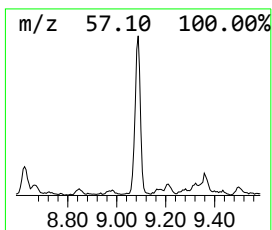
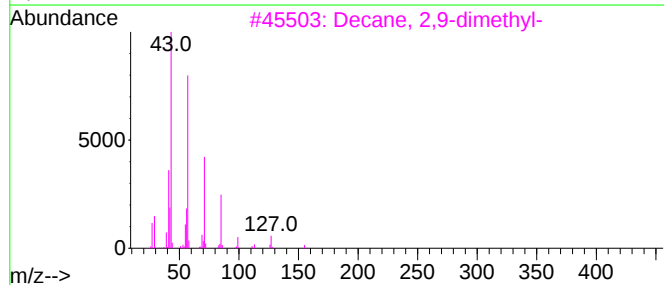
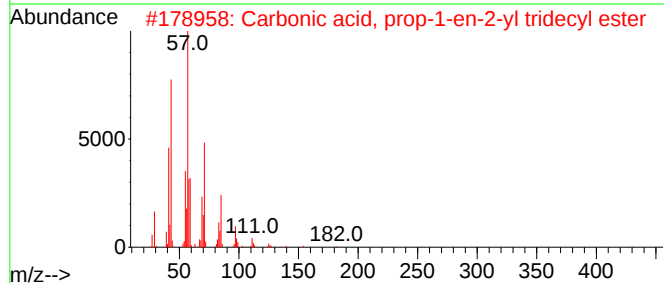
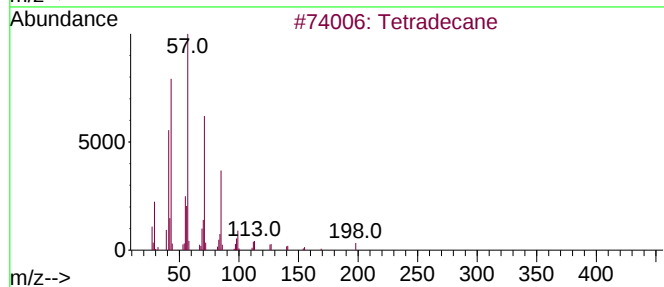
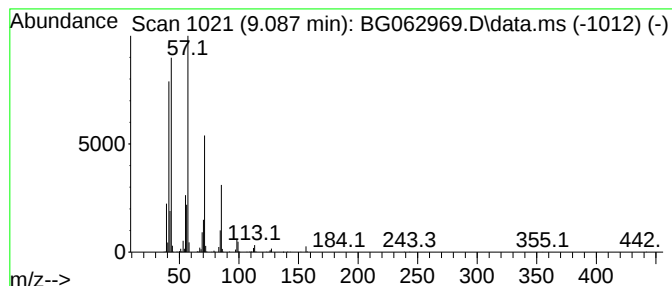
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.087	56.23 ng/ul	1512280	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	78
2			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	78
3			Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	64
4			Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	59
5			Undecane, 2-methyl-	170	C12H26	007045-71-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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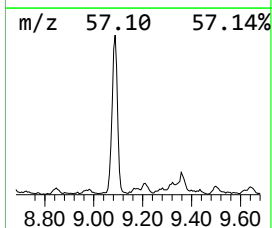
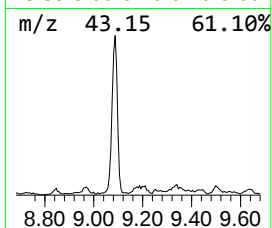
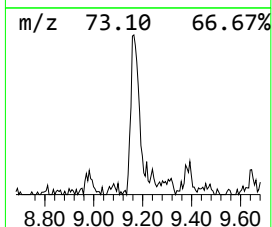
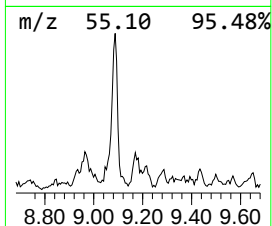
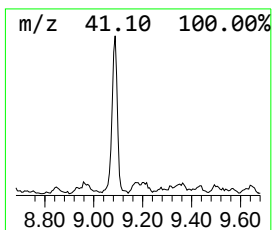
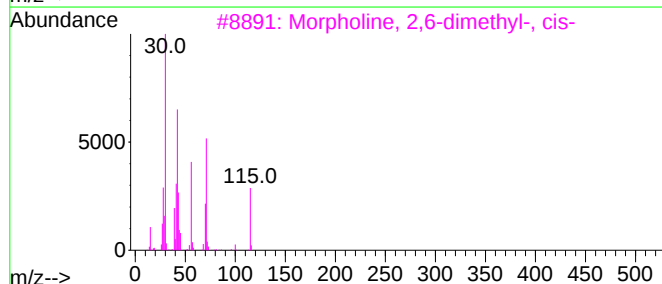
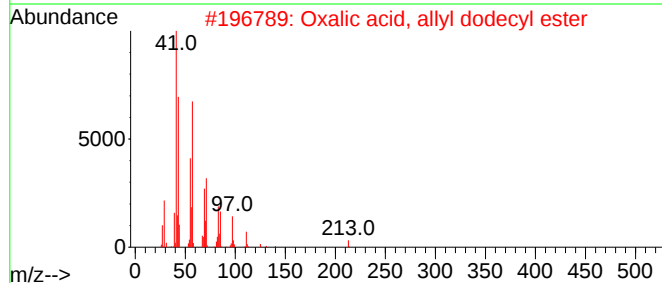
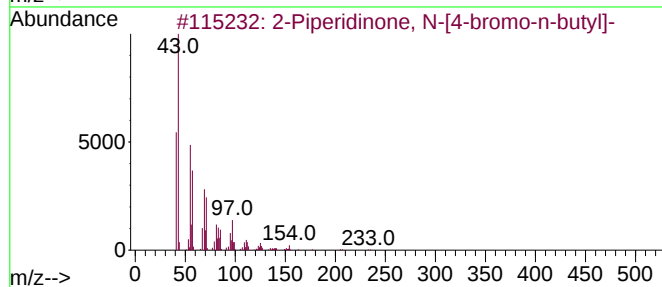
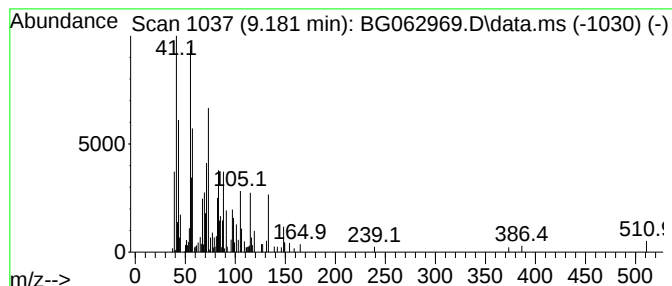
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-04 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.181	7.16 ng/ul	192681	1,4-Dichlorobenzene-d4	7.853

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Piperidinone, N-[4-bromo-n-but...	233	C9H16BrNO	195194-80-0	25		
2	Oxalic acid, allyl dodecyl ester	298	C17H30O4	1000309-24-0	22		
3	Morpholine, 2,6-dimethyl-, cis-	115	C6H13NO	006485-55-8	22		
4	Octadecyl octyl ether	382	C26H54O	1000406-38-7	22		
5	Oxalic acid, allyl hexadecyl ester	354	C21H38O4	1000309-24-4	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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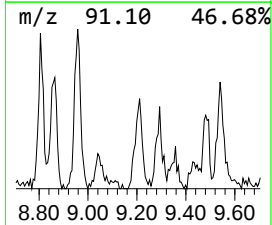
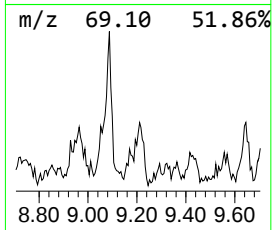
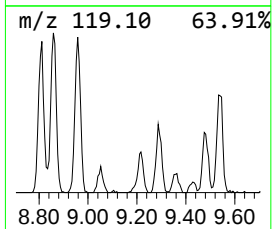
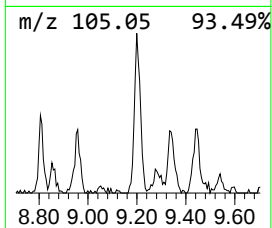
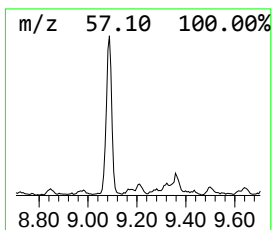
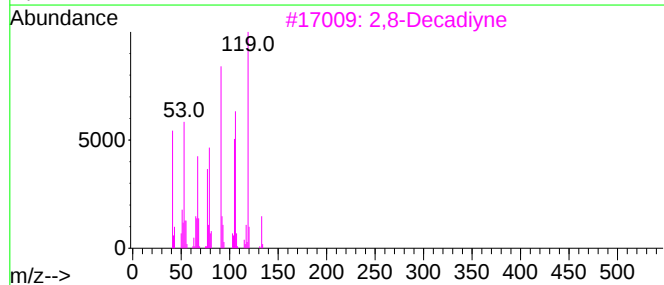
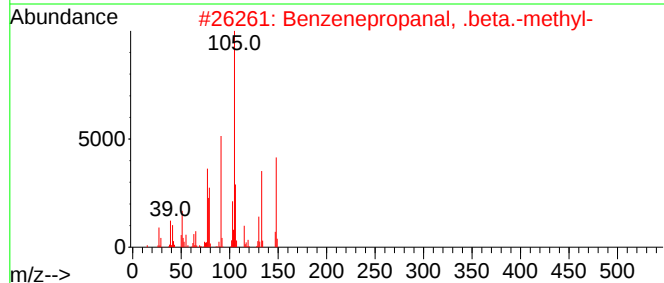
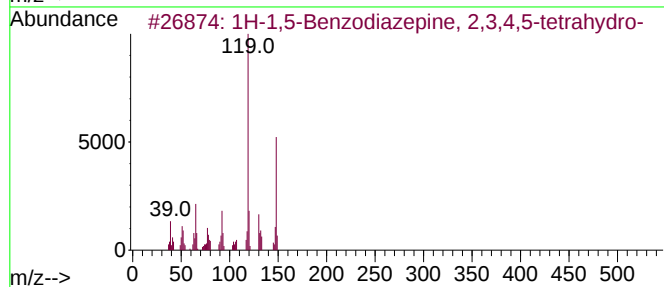
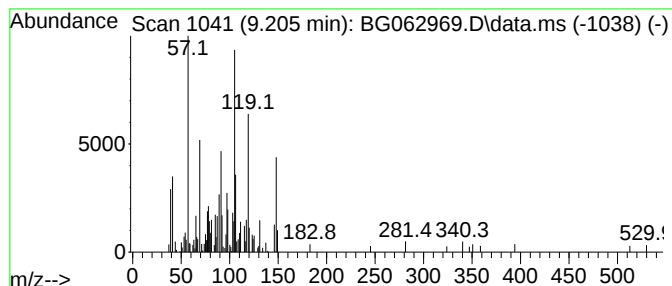
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-05 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.205	8.96 ng/ul	241044	1,4-Dichlorobenzene-d4	7.853

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-1,5-Benzodiazepine, 2,3,4,5-t...	148	C9H12N2	006516-89-8	38
2		Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	30
3		2,8-Decadiyne	134	C10H14	004116-93-2	25
4		Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	22
5		Benzene, (1,2-dimethylpropyl)-	148	C11H16	004481-30-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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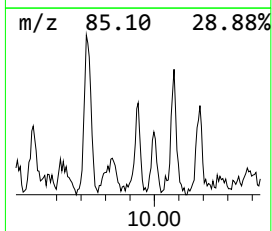
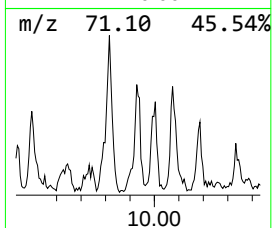
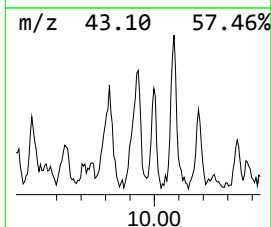
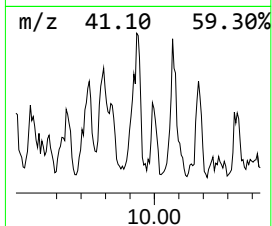
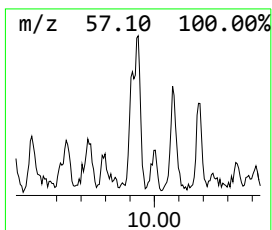
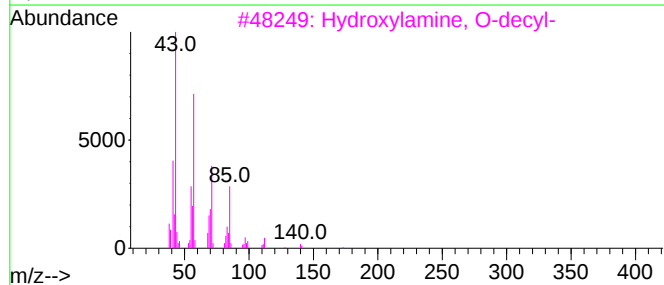
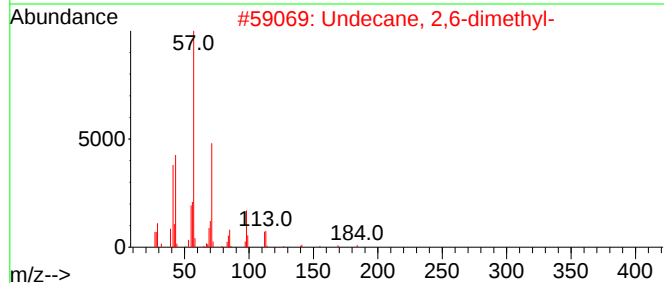
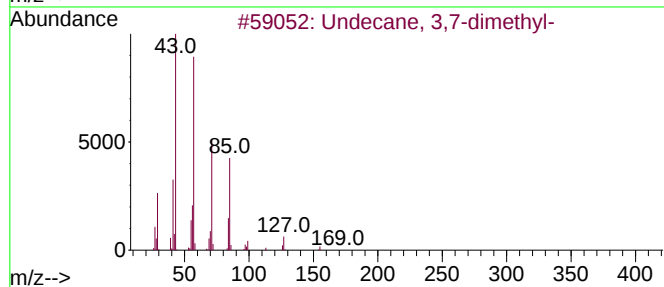
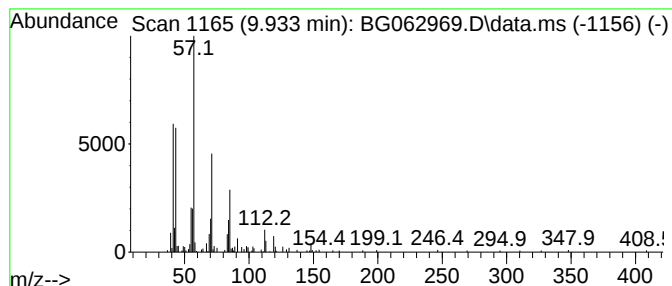
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 64

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.933	3.17 ng/ul	282768	Naphthalene-d8	10.644

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	64
2		Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	53
3		Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	53
4		Undecane, 5-methyl-	170	C12H26	001632-70-8	52
5		Tridecane, 6-methyl-	198	C14H30	013287-21-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

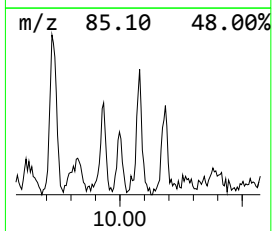
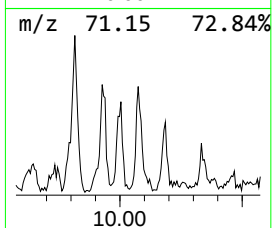
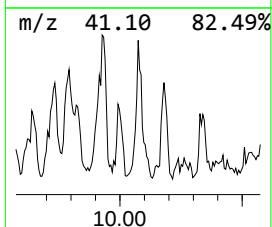
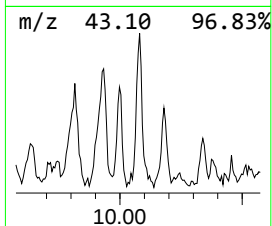
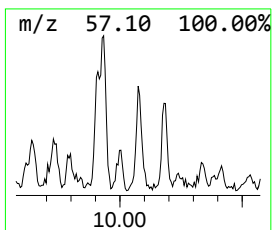
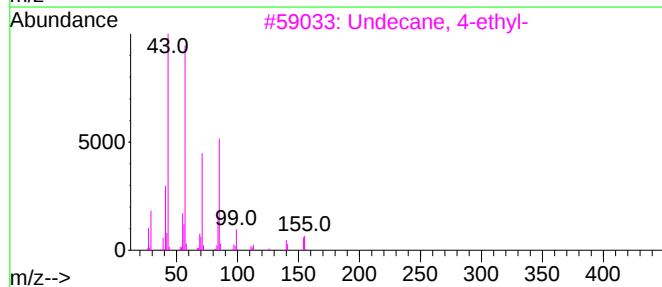
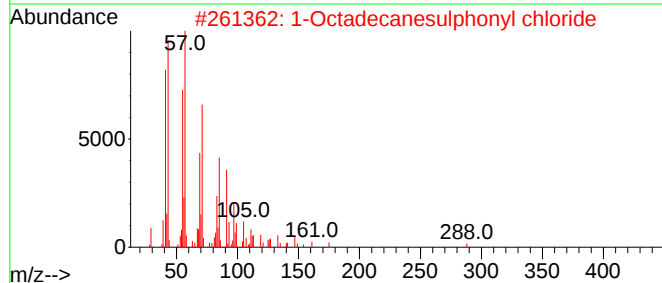
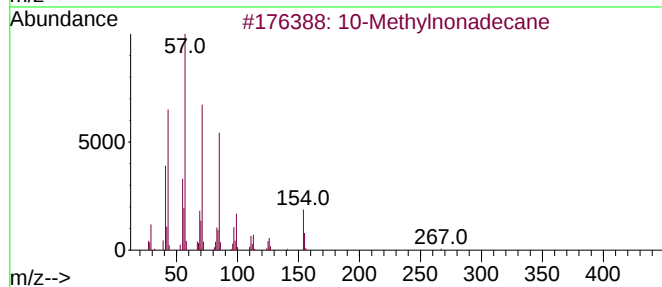
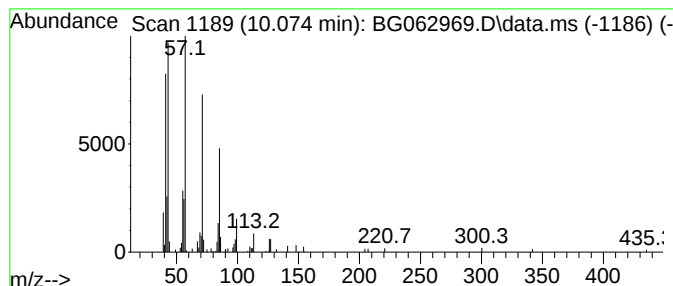
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.074	2.11 ng/ul	187841	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10-Methylnonadecane	282	C20H42	056862-62-5	78
2	1-Octadecanesulphonyl chloride	352	C18H37ClO2S	1000342-70-4	64
3	Undecane, 4-ethyl-	184	C13H28	017312-59-3	59
4	Hexadecane	226	C16H34	000544-76-3	59
5	Undecane, 3-methyl-	170	C12H26	001002-43-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
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Sample : P3898-20ME
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ALS Vial : 33 Sample Multiplier: 1

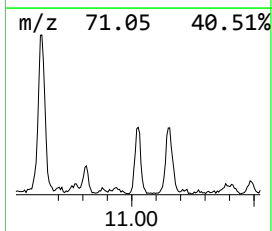
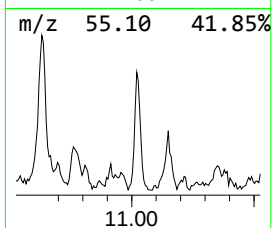
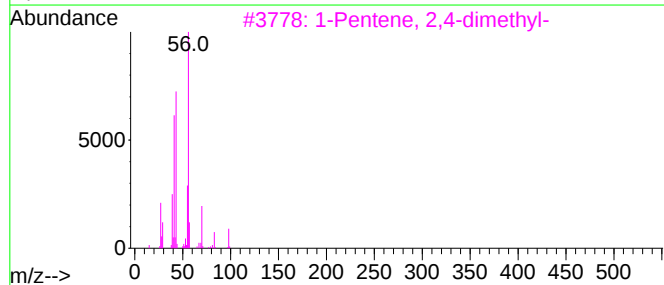
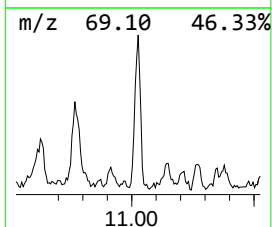
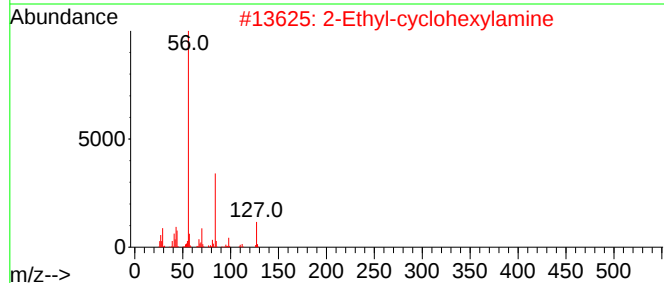
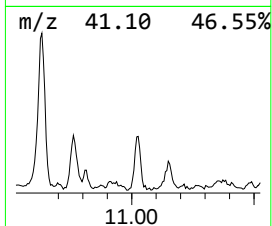
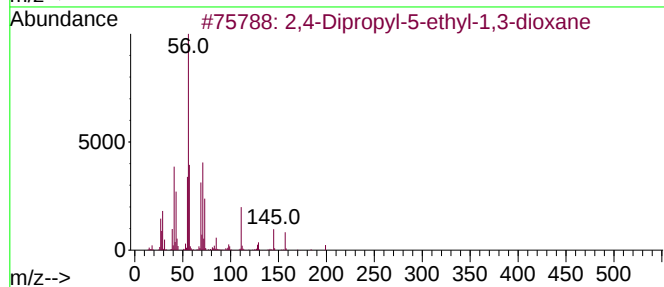
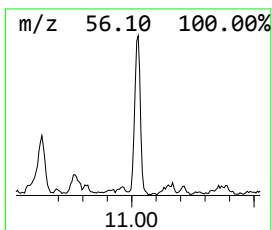
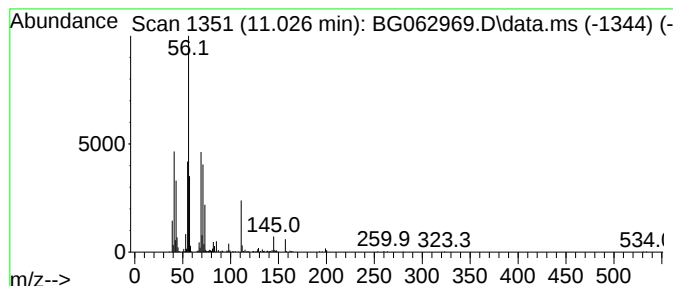
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 30 2,4-Dipropyl-5-ethyl-1,3-di... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.026	6.28 ng/ul	559565	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	64
2	2-Ethyl-cyclohexylamine	127	C8H17N	024216-90-8	35
3	1-Pentene, 2,4-dimethyl-	98	C7H14	002213-32-3	35
4	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	35
5	2-Pyrrolidinone, 1-ethenyl-	111	C6H9NO	000088-12-0	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

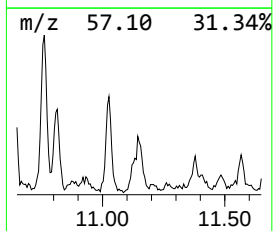
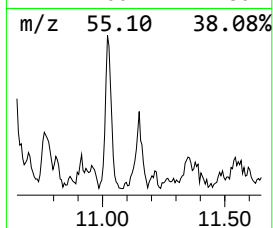
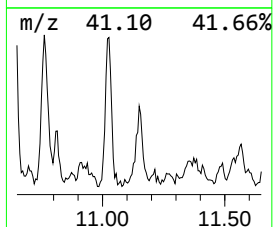
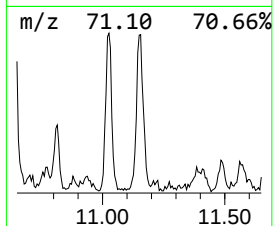
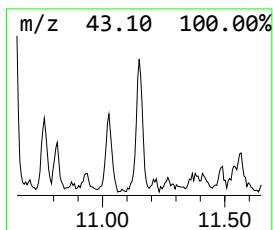
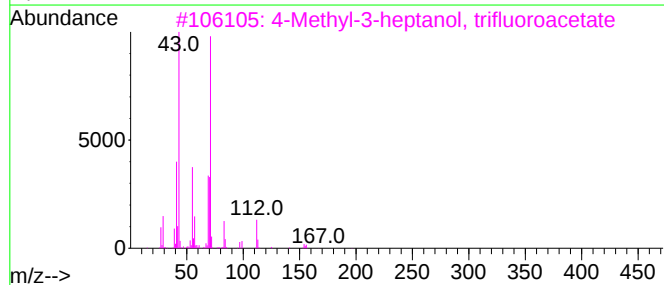
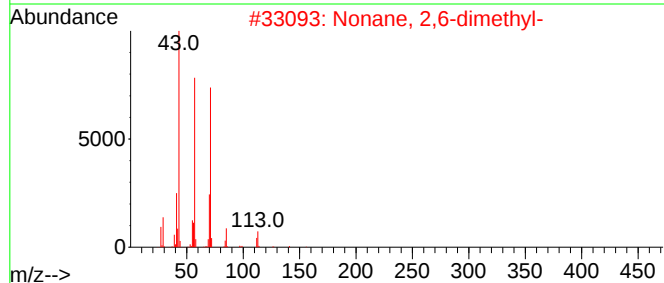
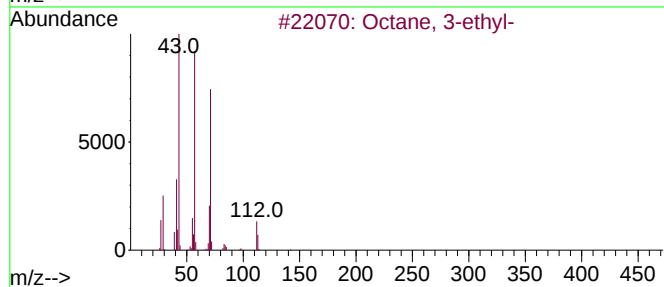
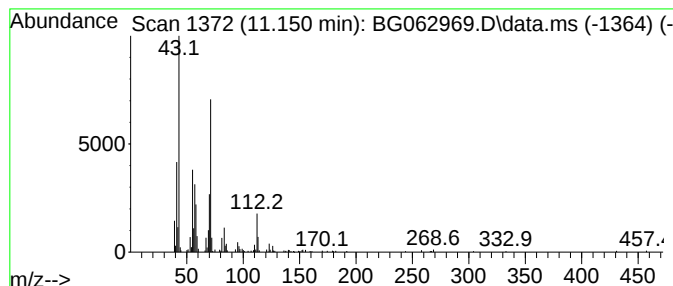
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BNA_G
ClientSampleId :
DDC59ME

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 31 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.150	3.89 ng/ul	347145	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-ethyl-	142	C10H22	005881-17-4	70
2	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	50
3	4-Methyl-3-heptanol, trifluoroac...	226	C10H17F3O2	1000365-51-8	50
4	Sulfurous acid, octadecyl 2-pent...	404	C23H48O3S	1000309-16-6	43
5	Butyric acid, m-fluorophenyl ester	182	C10H11FO2	029052-04-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

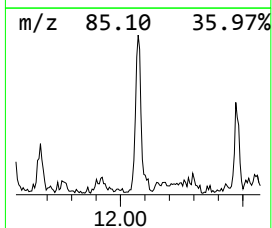
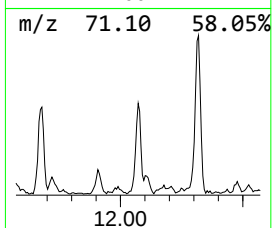
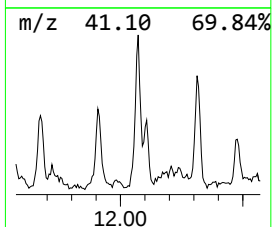
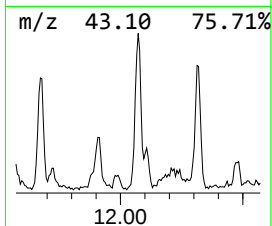
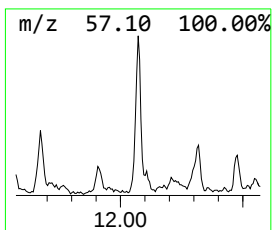
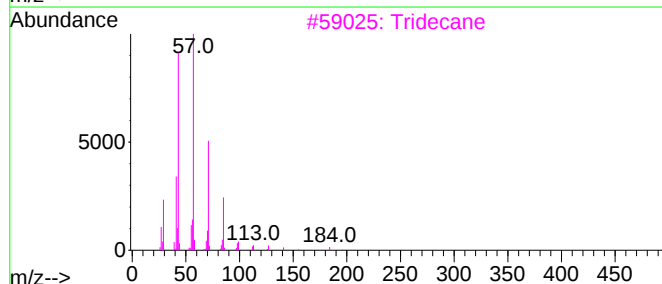
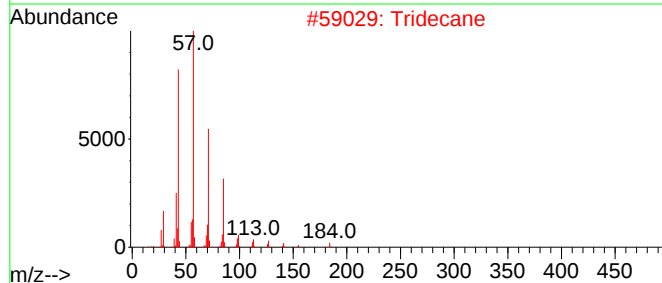
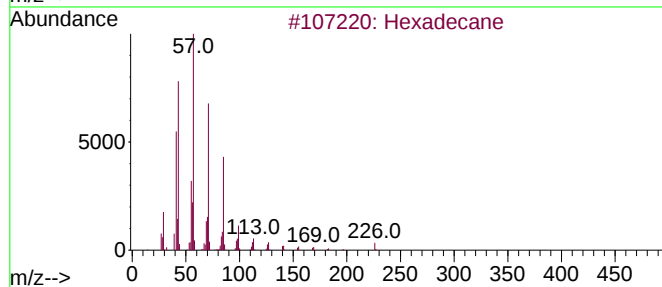
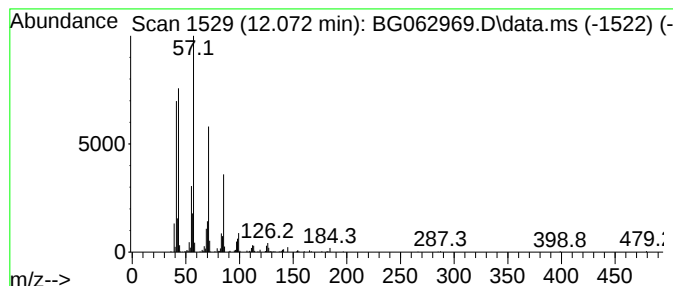
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.072	6.06 ng/ul	540337	Naphthalene-d8	10.644	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	87
2	Tridecane	184	C13H28	000629-50-5	87
3	Tridecane	184	C13H28	000629-50-5	87
4	Nonadecane	268	C19H40	000629-92-5	86
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

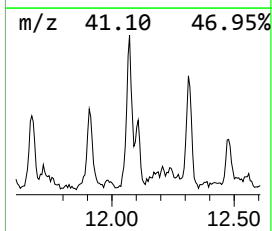
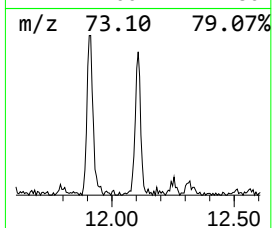
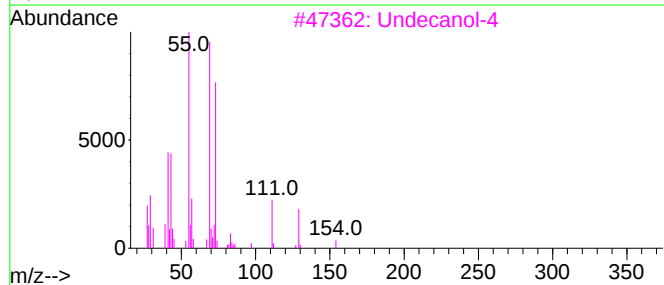
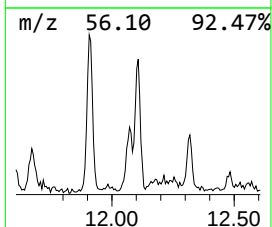
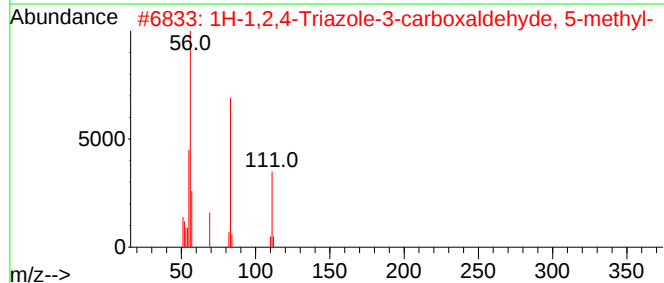
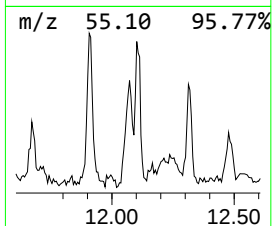
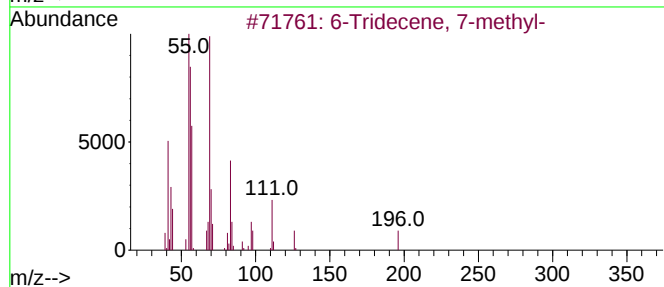
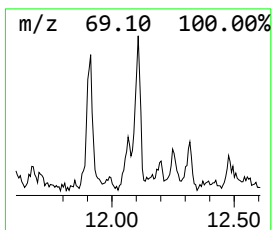
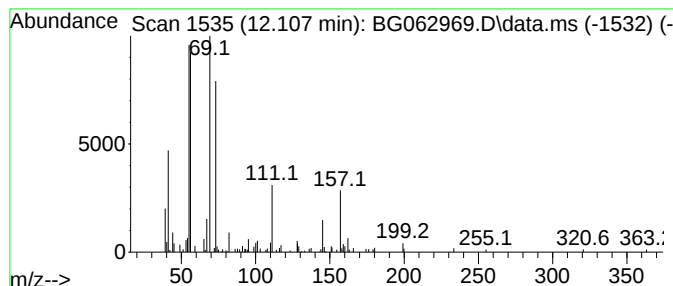
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.107	2.71 ng/ul	241765	Naphthalene-d8	10.644

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Tridecene, 7-methyl-			196	C14H28	024949-42-6	38
2	1H-1,2,4-Triazole-3-carboxaldehy...			111	C4H5N3O	056804-98-9	27
3	Undecanol-4			172	C11H24O	004272-06-4	25
4	6-Tridecanol			200	C13H28O	005770-03-6	23
5	2,3,4-Trimethyl-hex-3-enal			140	C9H16O	1000193-72-9	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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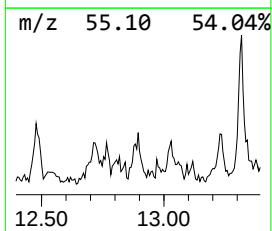
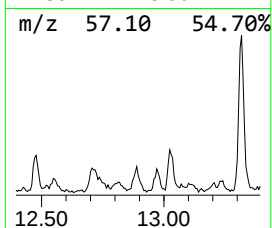
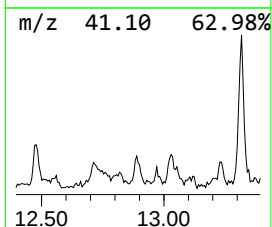
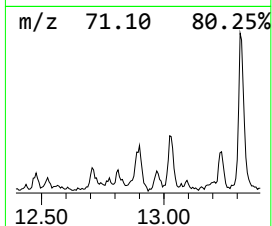
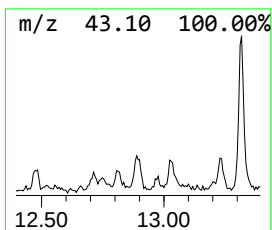
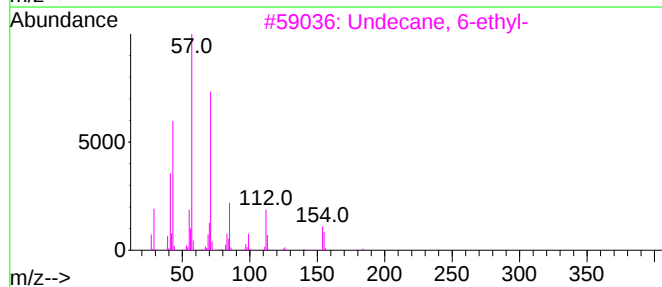
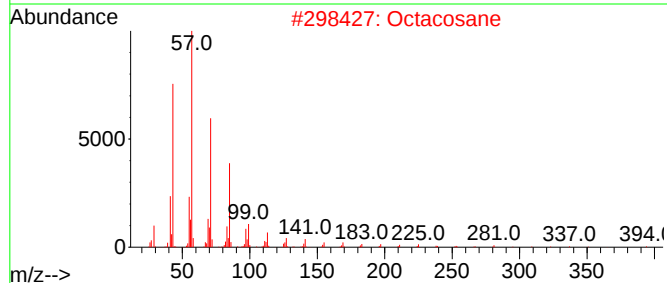
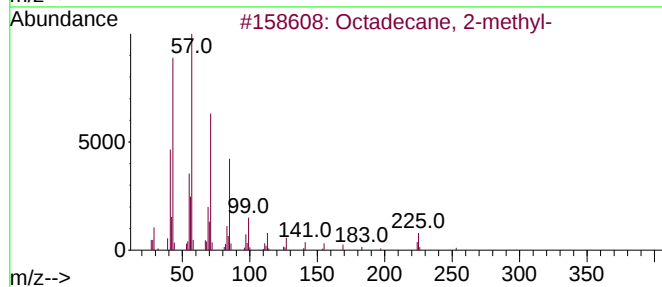
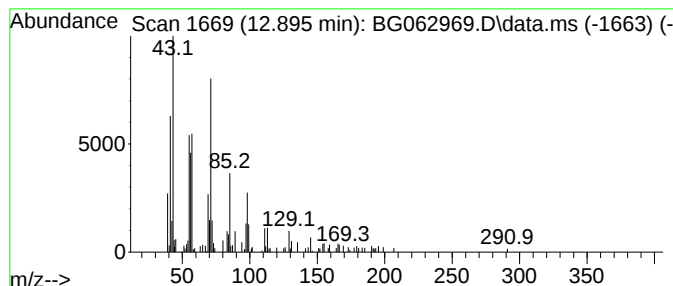
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.895	4.86 ng/ul	145934	Acenaphthene-d10	14.487

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane, 2-methyl-	268	C19H40	001560-88-9	50
2		Octacosane	394	C28H58	000630-02-4	50
3		Undecane, 6-ethyl-	184	C13H28	017312-60-6	47
4		Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	47
5		Heptacosane	380	C27H56	000593-49-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

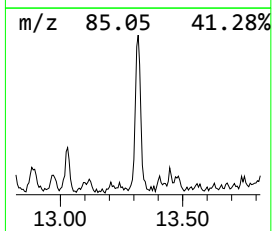
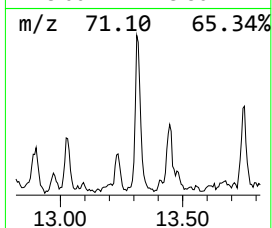
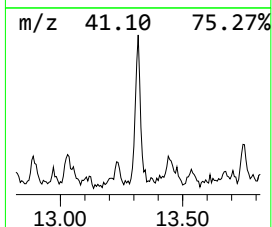
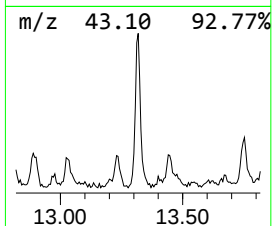
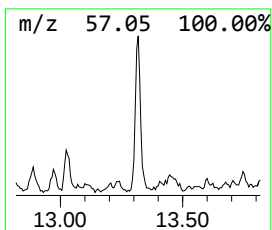
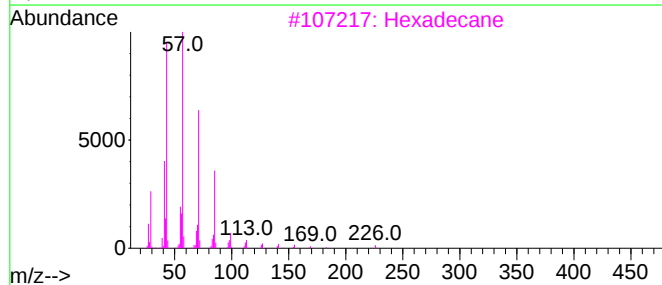
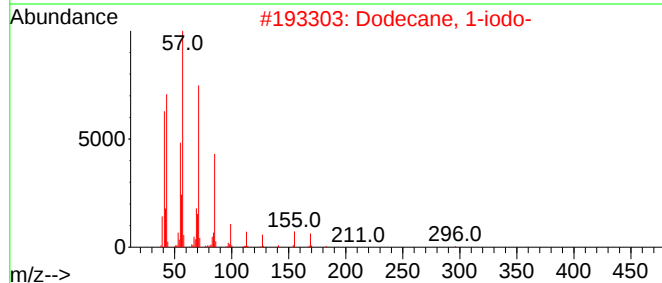
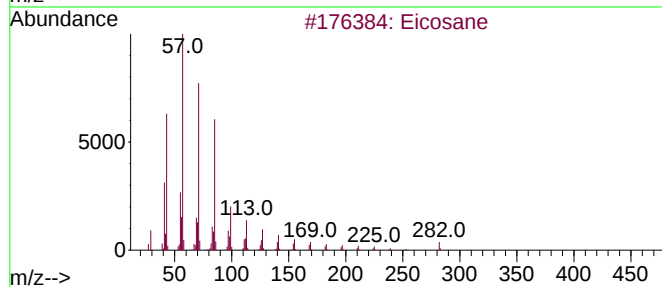
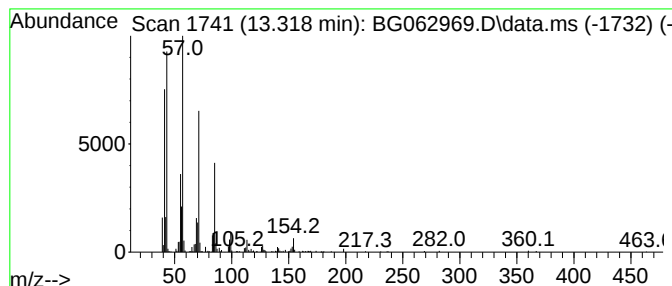
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
13.318	19.33 ng/ul	580623	Acenaphthene-d10		14.487	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	90
2	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	87
3	Hexadecane		226	C16H34	000544-76-3	87
4	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	87
5	Tridecane		184	C13H28	000629-50-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

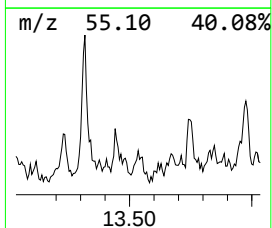
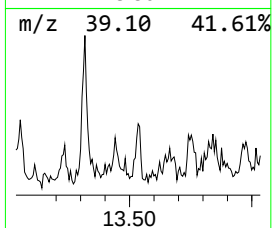
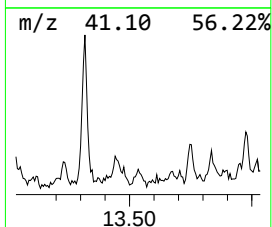
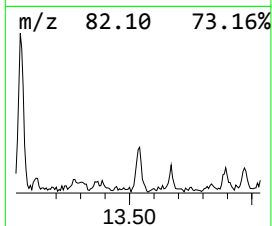
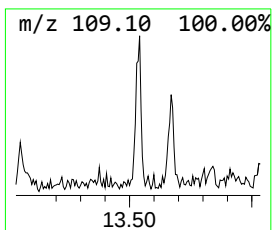
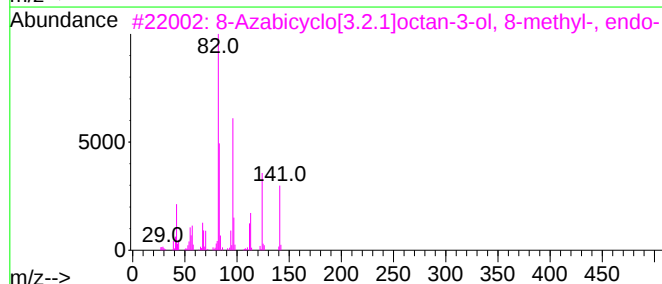
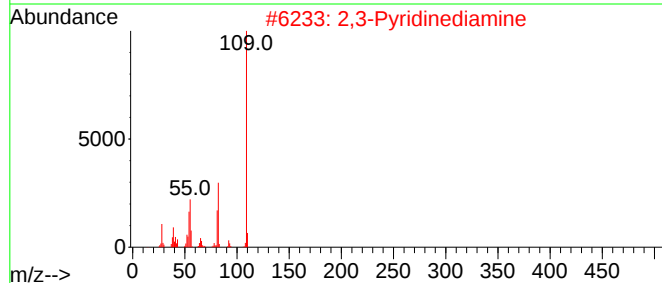
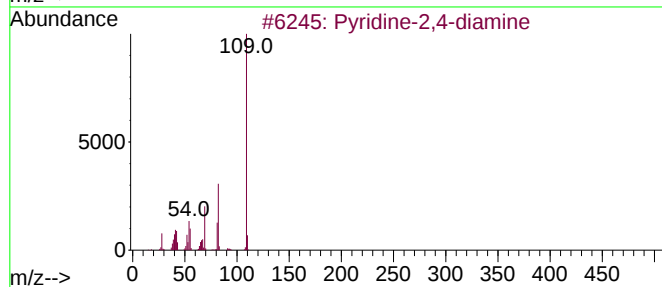
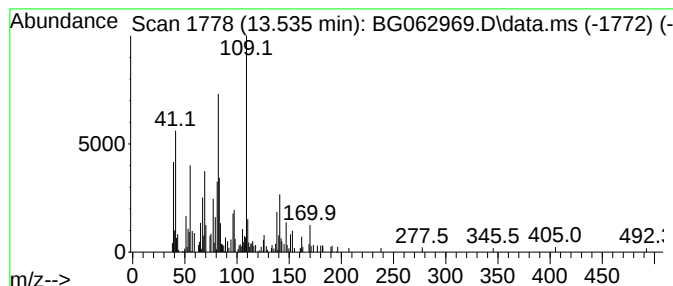
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Pyridine-2,4-diamine Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.535	6.91 ng/ul	207502	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyridine-2,4-diamine	109	C5H7N3	000461-88-1	50
2			2,3-Pyridinediamine	109	C5H7N3	000452-58-4	50
3			8-Azabicyclo[3.2.1]octan-3-ol, 8...	141	C8H15NO	000120-29-6	35
4			4-Amino-2-methylpyrimidine	109	C5H7N3	000074-69-1	27
5			8-Azabicyclo[3.2.1]octan-3-ol, 8...	141	C8H15NO	000135-97-7	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

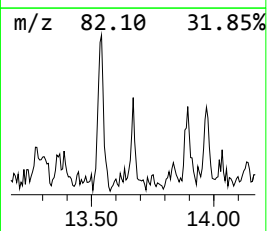
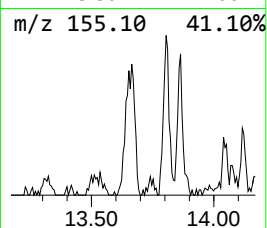
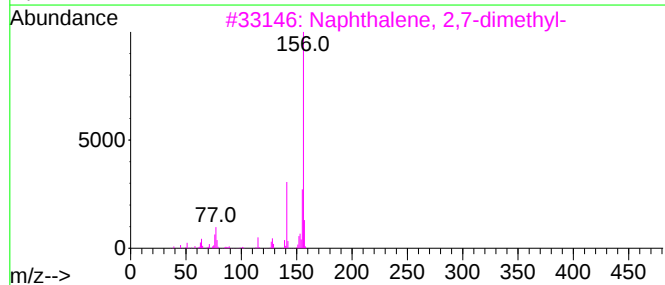
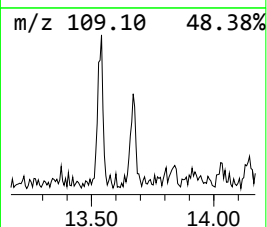
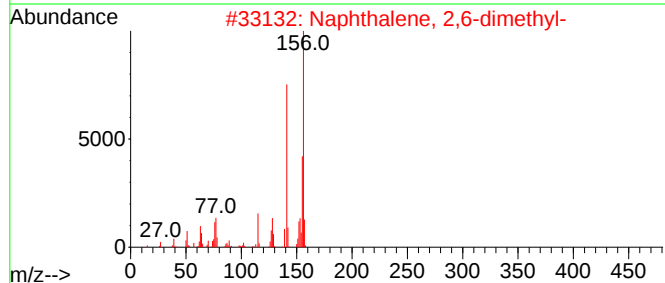
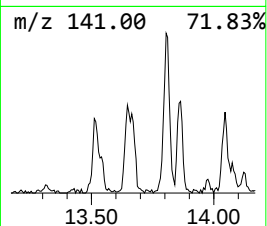
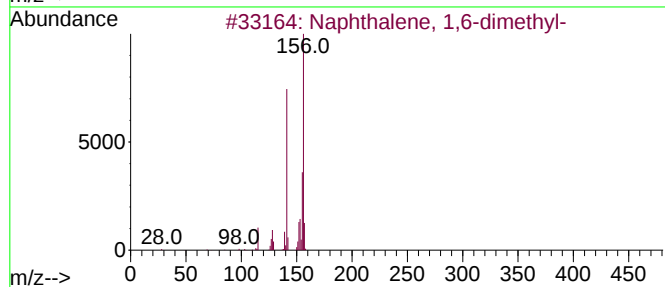
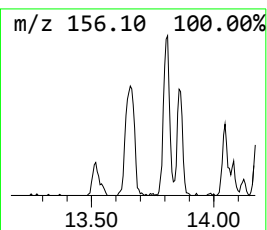
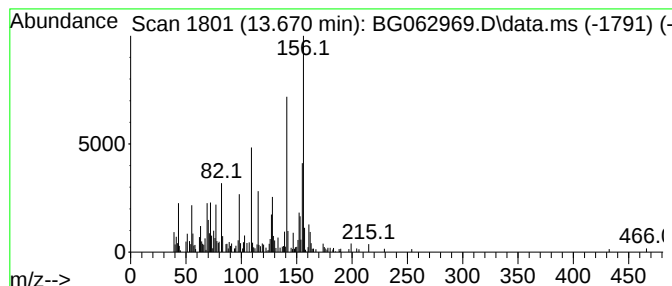
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 Naphthalene, 1,6-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.670	9.61 ng/ul	288788	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	83
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	78
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	78
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	64
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

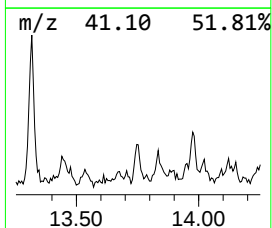
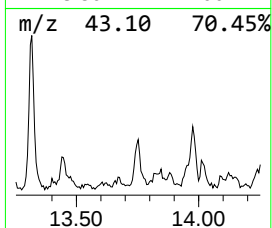
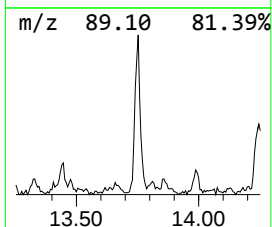
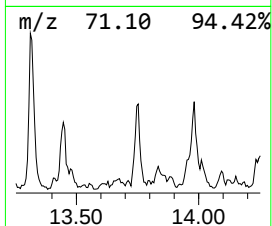
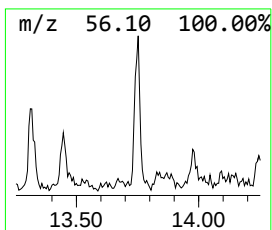
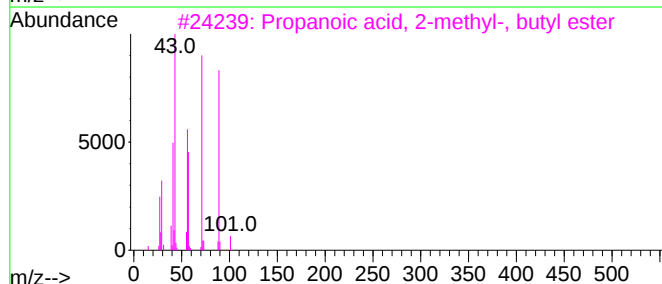
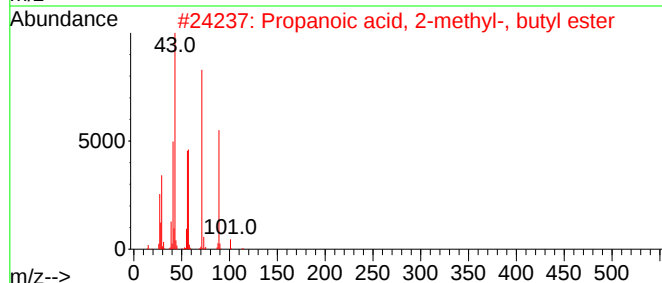
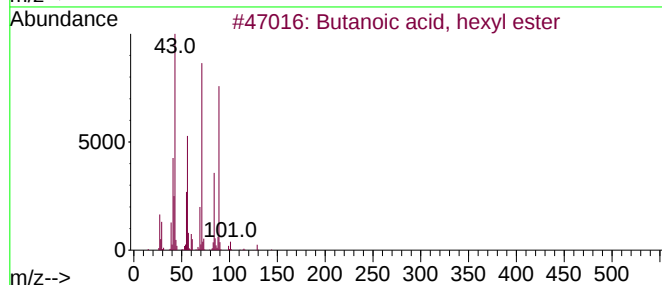
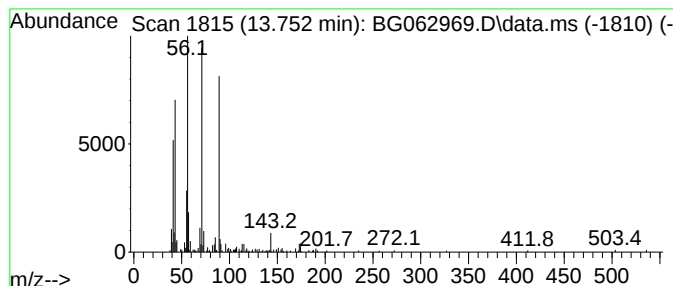
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 Butanoic acid, hexyl ester Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.752	6.48 ng/ul	194762	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	72
2			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	72
3			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	72
4			Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	56
5			Propanoic acid, 2-methyl-, butyl...	144	C8H16O2	000097-87-0	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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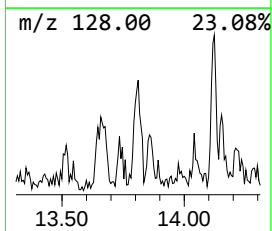
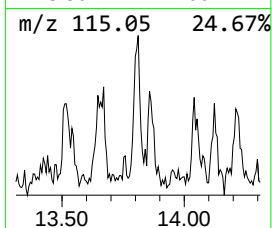
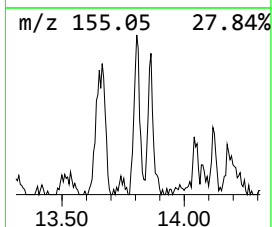
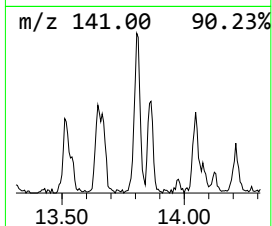
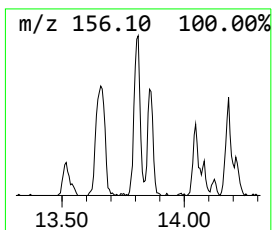
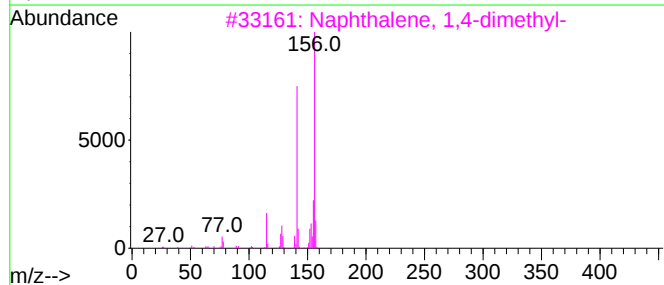
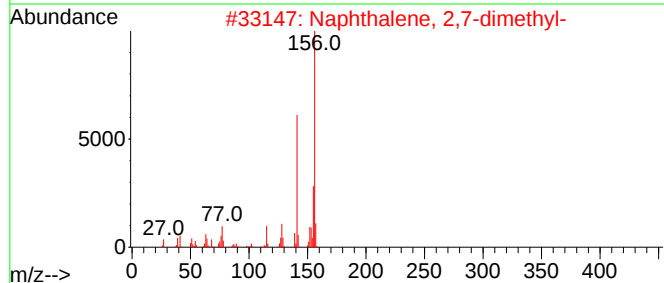
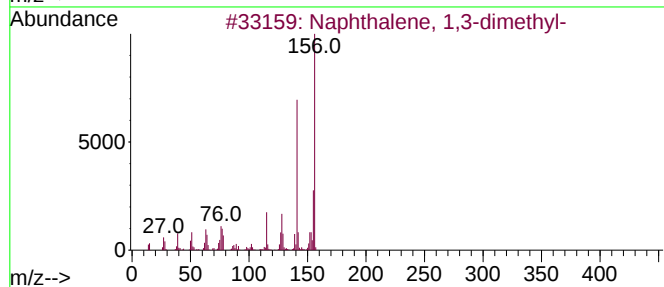
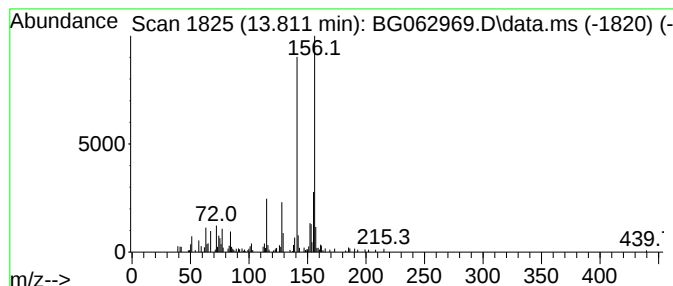
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Naphthalene, 1,3-dimethyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.811	5.39 ng/ul	161762	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	91
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	91
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	90
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	89
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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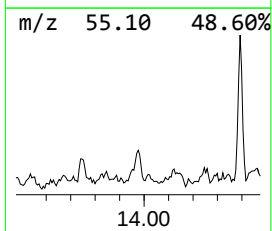
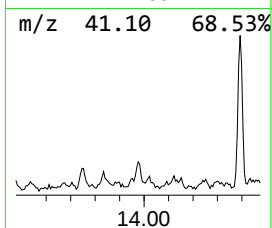
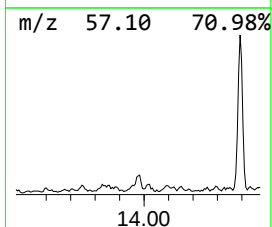
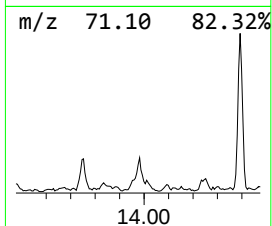
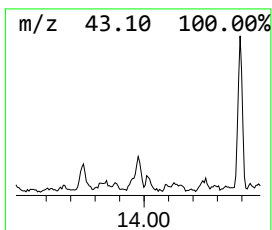
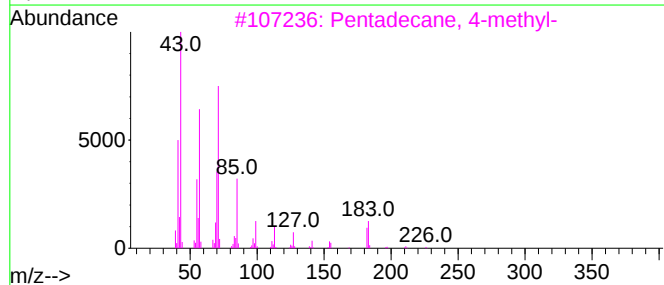
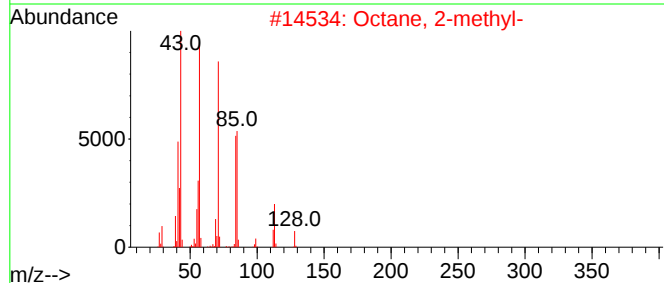
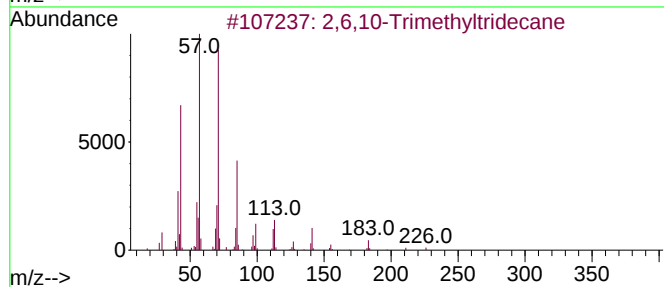
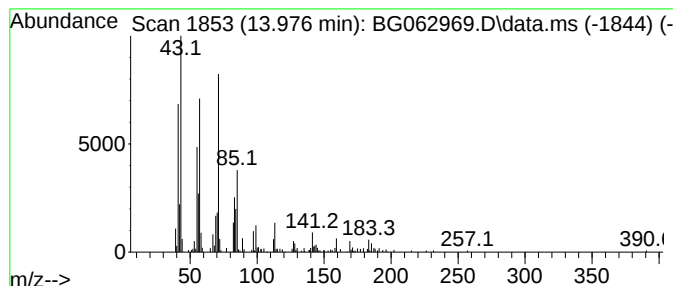
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD		R.T.
13.976	9.13 ng/ul	274372	Acenaphthene-d10		14.487

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Trimethyltridecane	226	C16H34		003891-99-4	92
2	Octane, 2-methyl-	128	C9H20		003221-61-2	70
3	Pentadecane, 4-methyl-	226	C16H34		002801-87-8	70
4	Tetradecane, 4,11-dimethyl-	226	C16H34		055045-12-0	62
5	Decane, 3,6-dimethyl-	170	C12H26		017312-53-7	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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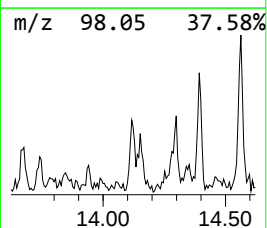
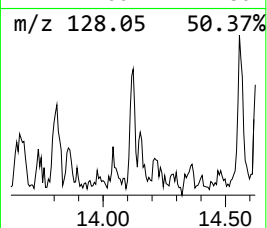
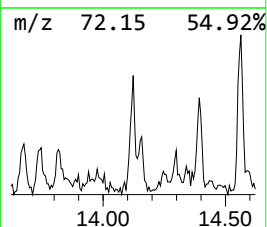
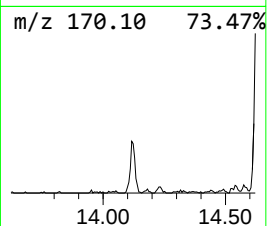
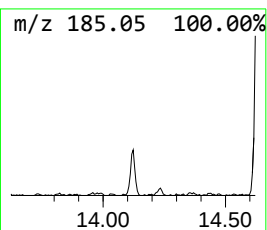
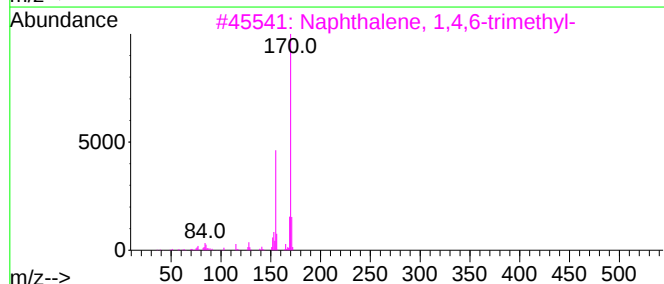
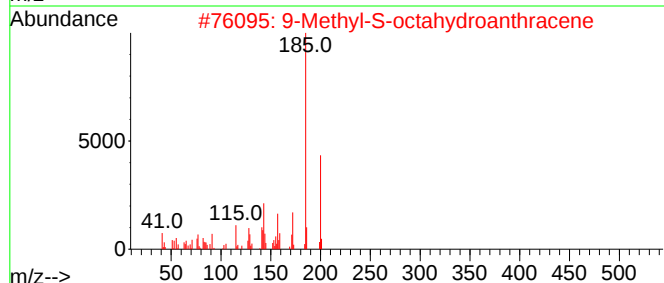
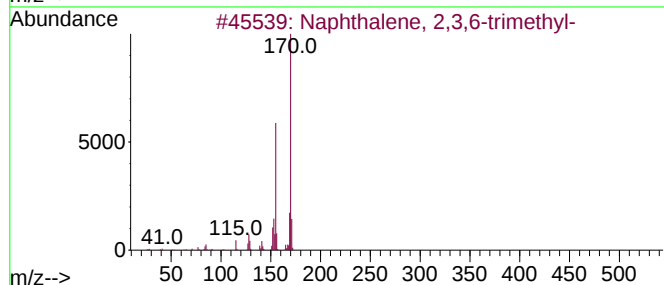
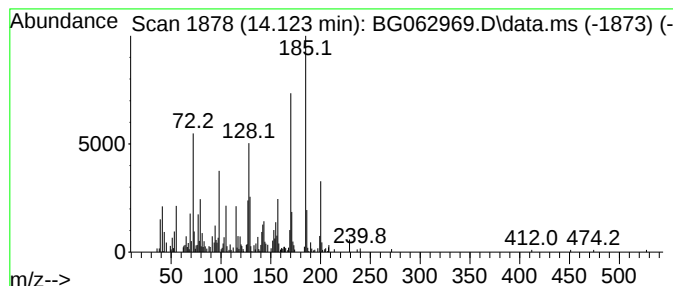
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TIC Integration Parameters: LSCINT.P

Peak Number 44 unknown-06

Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.123	7.02 ng/ul	210827	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	42
2			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	38
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	30
4			Silane, trimethyl-2-naphthalenyl-	200	C13H16Si	018052-85-2	25
5			3,4-Dimethoxythiophenol	170	C8H10O2S	000700-96-9	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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Sample : P3898-20ME
Misc :
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Instrument :
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ClientSampleId :
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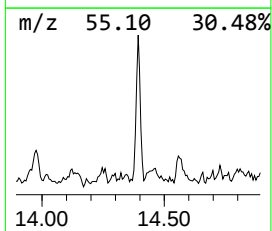
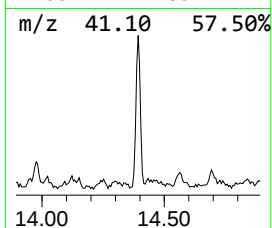
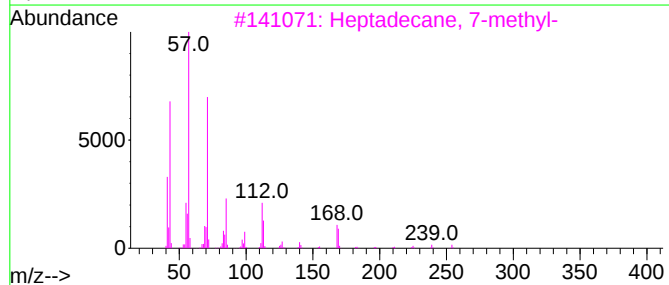
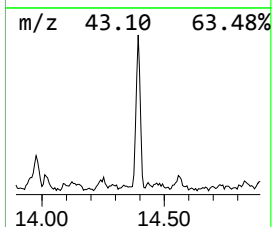
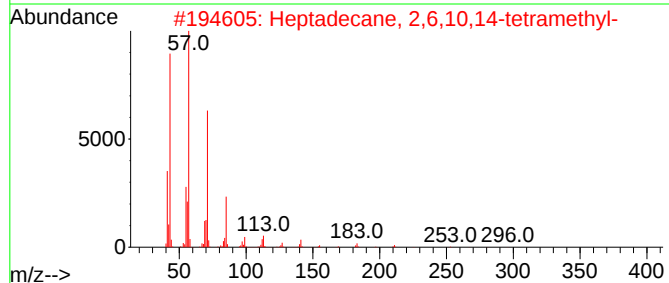
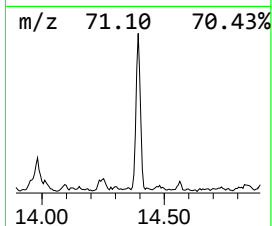
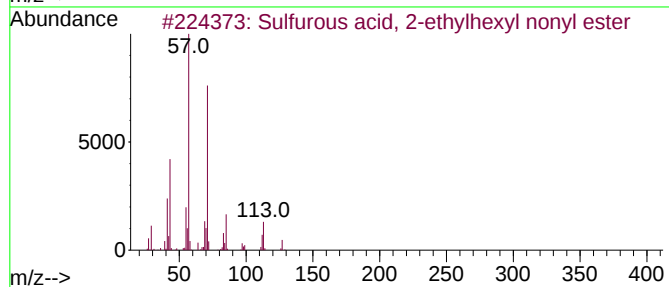
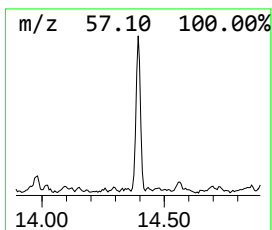
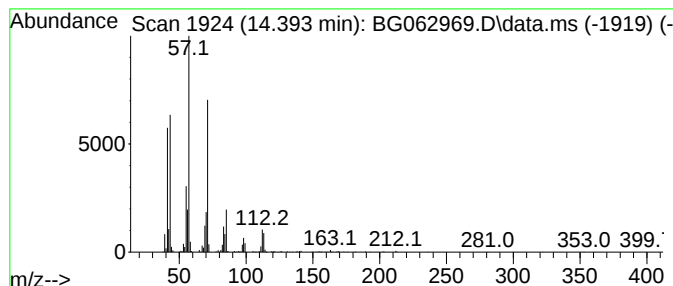
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 Sulfurous acid, 2-ethylhexy... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.393	31.27 ng/ul	939274	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72
2			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	72
3			Heptadecane, 7-methyl-	254	C18H38	020959-33-5	72
4			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72
5			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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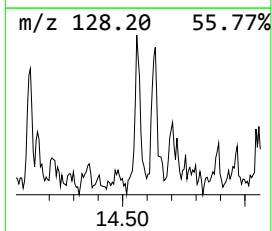
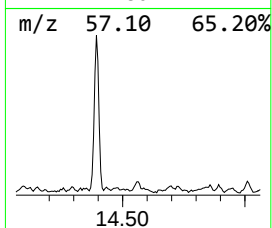
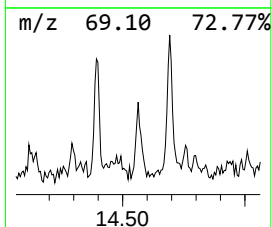
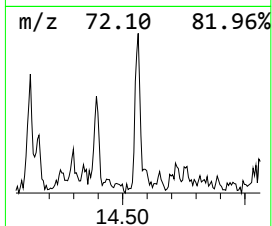
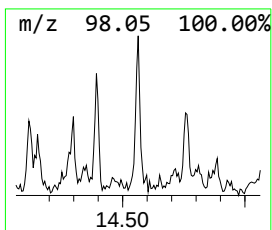
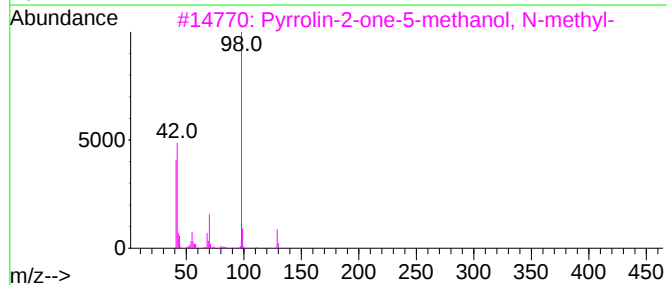
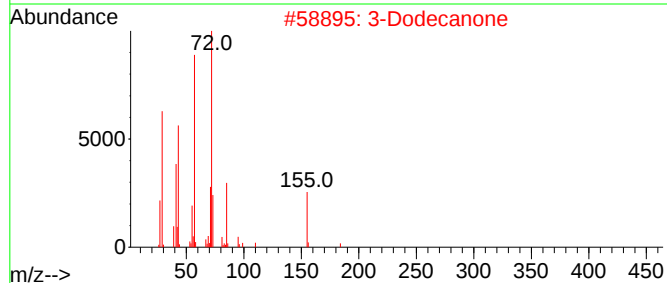
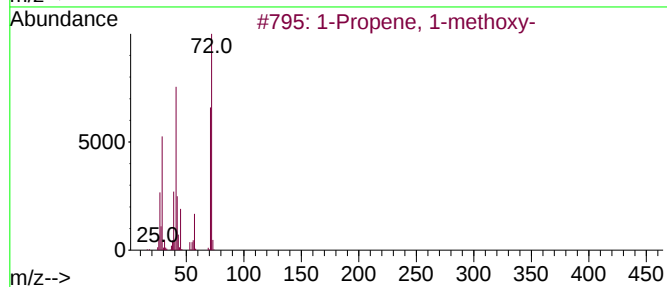
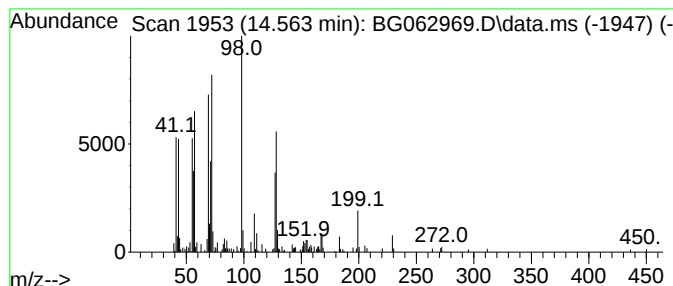
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Quant Title : SVOA CALIBRATION

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

Peak Number 46 unknown-07 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.563	7.58 ng/ul	227753	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propene, 1-methoxy-	72	C4H8O	007319-16-6	27
2			3-Dodecanone	184	C12H24O	001534-27-6	14
3			Pyrrolin-2-one-5-methanol, N-met...	129	C6H11NO2	1000125-98-2	14
4			2-Propanol, 1-octyloxy-(1-piperi...	271	C16H33NO2	096450-89-4	14
5			3-Piperidino-1,2-propanediol	159	C8H17NO2	004847-93-2	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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Misc :
ALS Vial : 33 Sample Multiplier: 1

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ClientSampleId :
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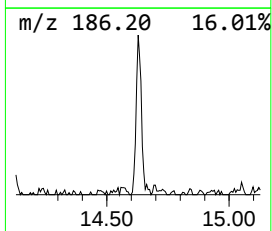
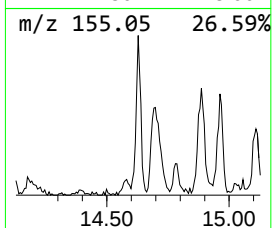
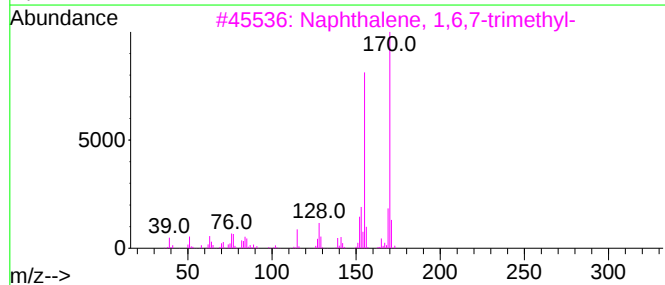
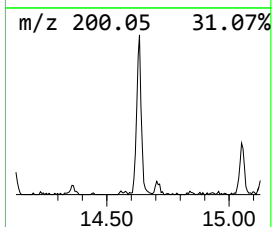
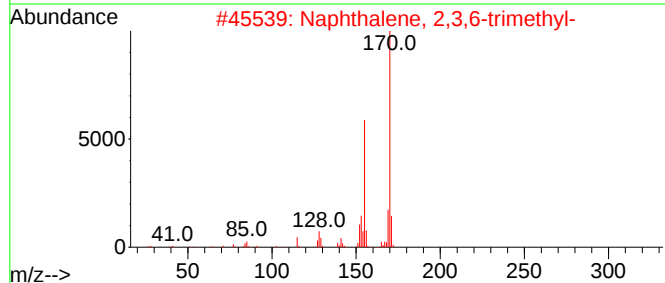
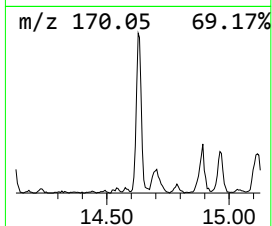
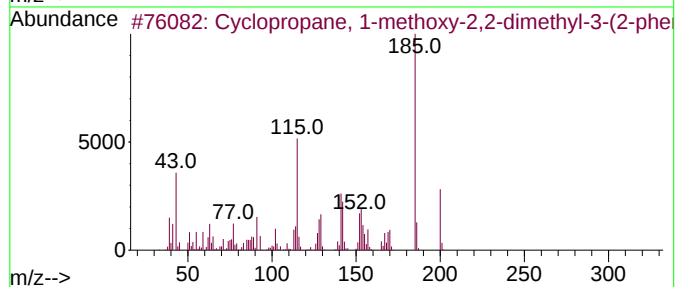
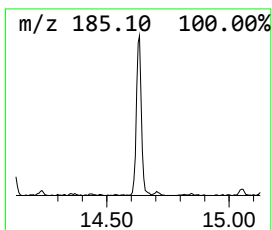
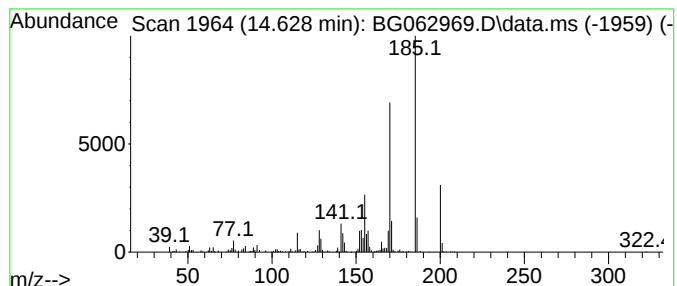
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 Cyclopropane, 1-methoxy-2,2... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.628	16.61 ng/ul	498979	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1-methoxy-2,2-dime...	200	C14H16O	1000271-83-0	64
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	49
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	47
4			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	47
5			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

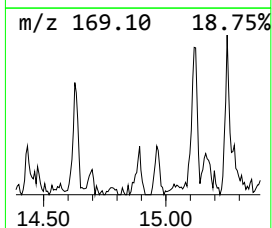
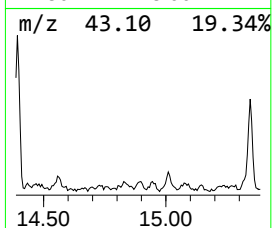
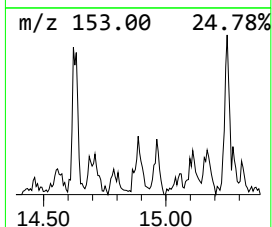
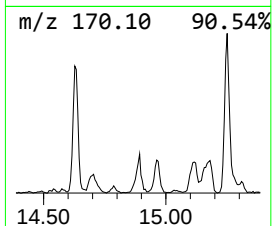
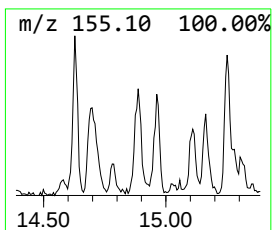
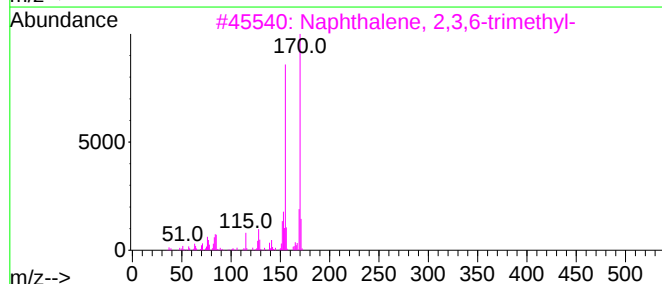
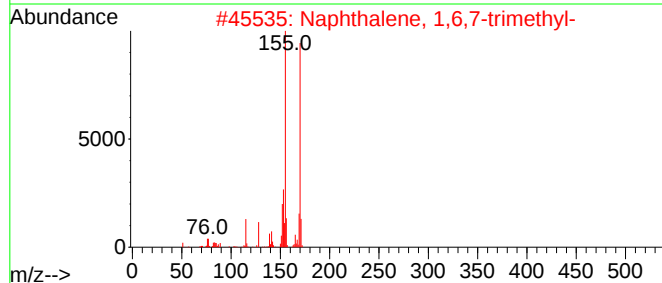
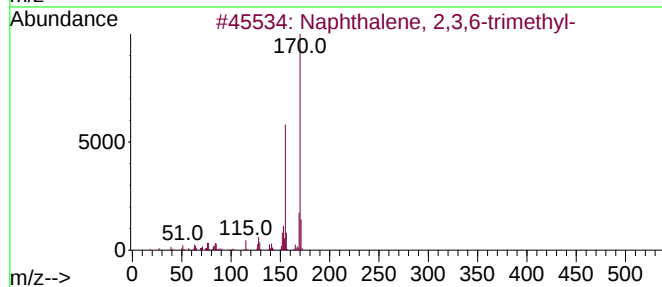
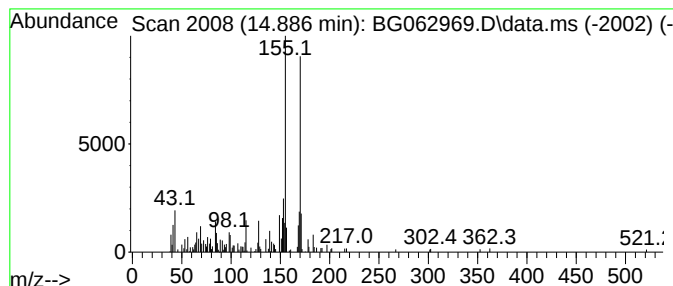
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 Naphthalene, 2,3,6-trimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.886	7.41 ng/ul	222657	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
5			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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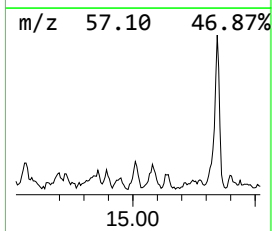
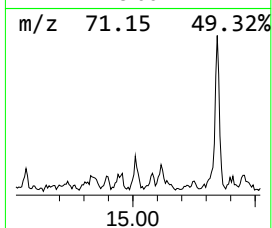
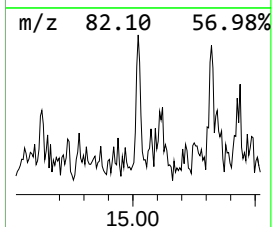
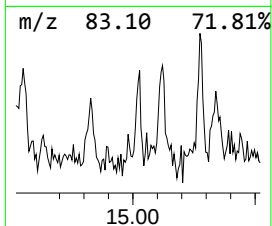
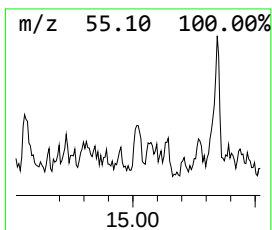
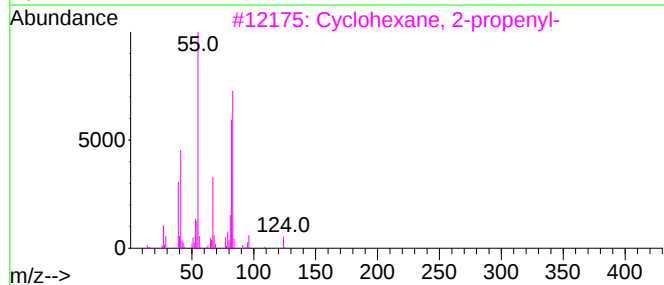
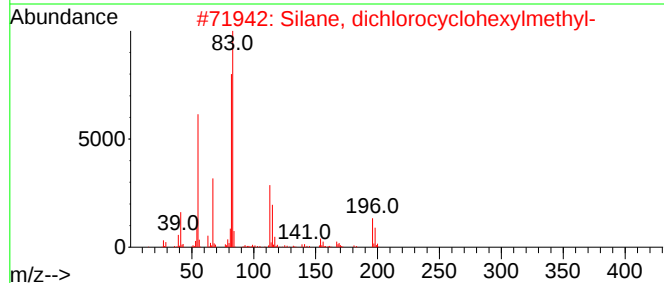
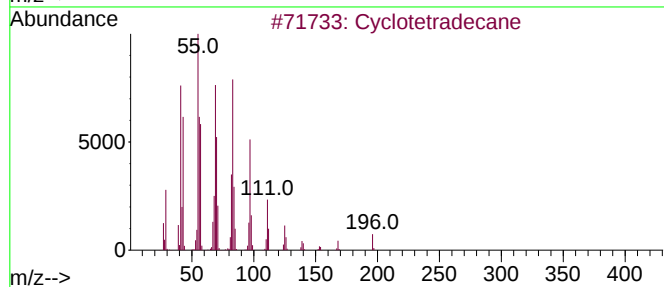
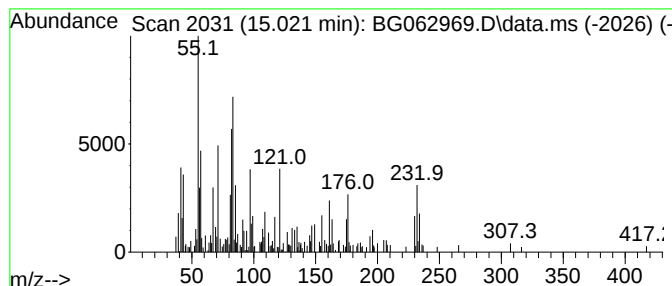
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 (DEL) Alkane: Cyclic15.022 Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.022	5.18 ng/ul	155646	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetradecane	196	C14H28	000295-17-0	43
2			Silane, dichlorocyclohexylmethyl-	196	C7H14Cl2Si	005578-42-7	30
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	30
4			Cyclohexane, octyl-	196	C14H28	001795-15-9	30
5			Cyclohexane, octyl-	196	C14H28	001795-15-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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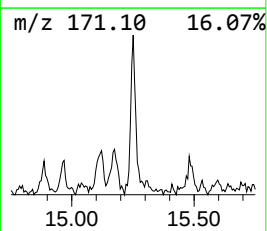
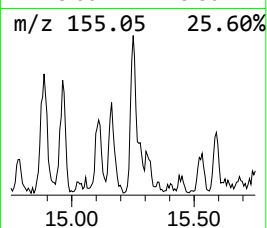
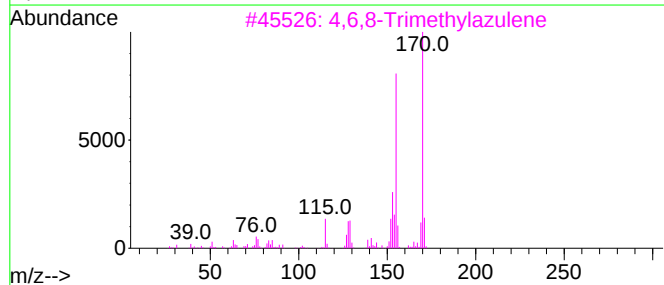
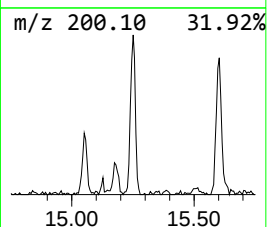
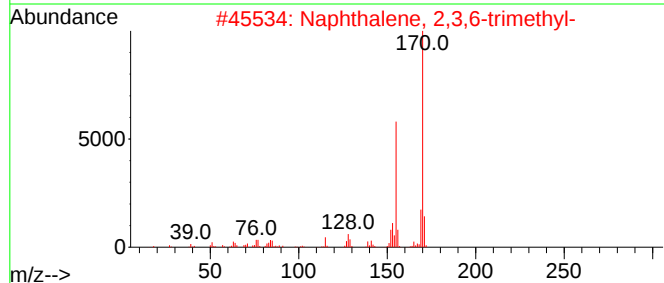
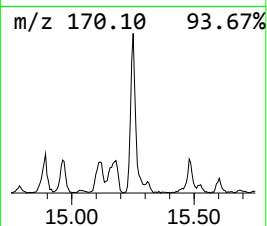
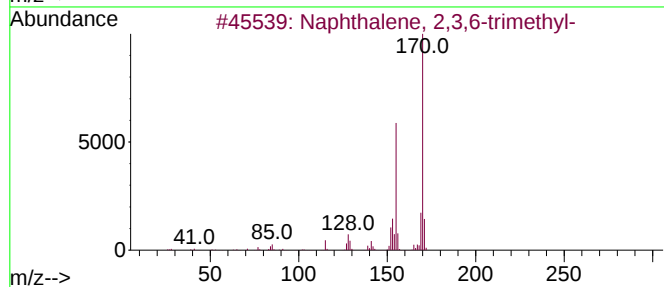
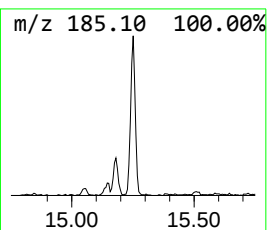
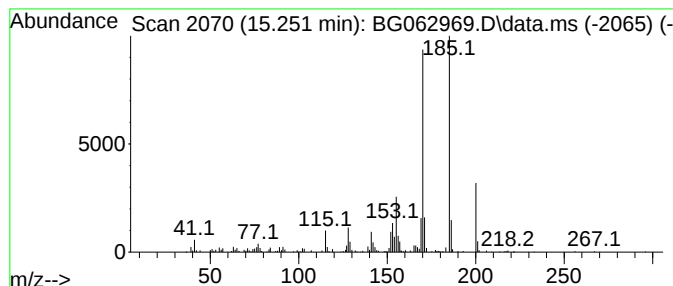
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.251	18.16 ng/ul	545521	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	86
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	64
3			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50
5			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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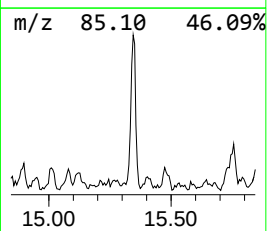
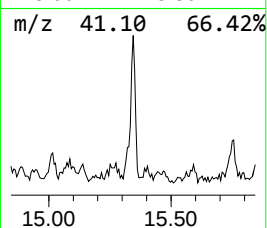
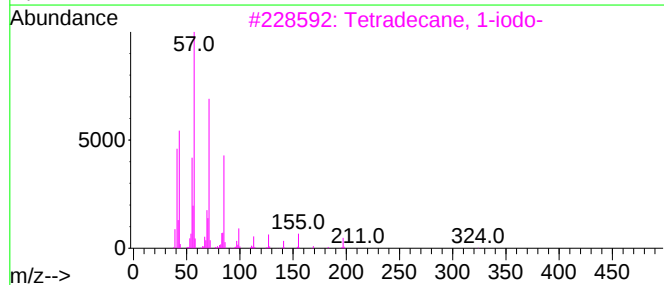
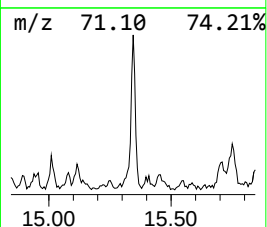
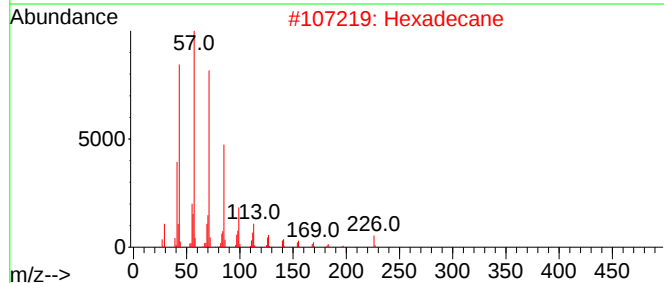
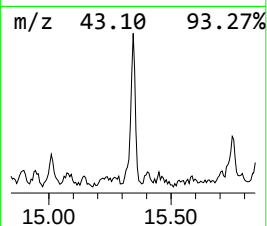
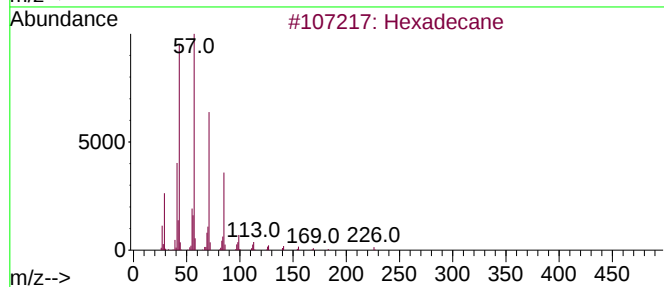
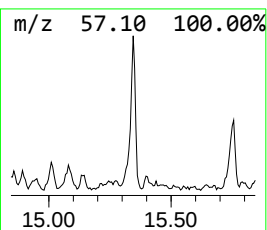
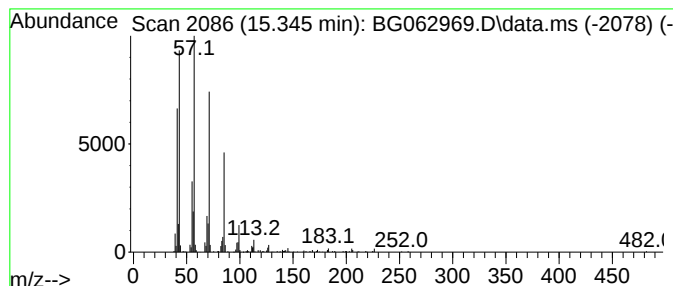
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Quant Title : SVOA CALIBRATION

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

Peak Number 51 Tetradecane, 1-iodo- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.345	19.42 ng/ul	583395	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	96
2			Hexadecane	226	C16H34	000544-76-3	94
3			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	91
4			Hexadecane	226	C16H34	000544-76-3	91
5			Tetradecane	198	C14H30	000629-59-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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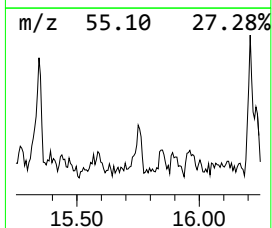
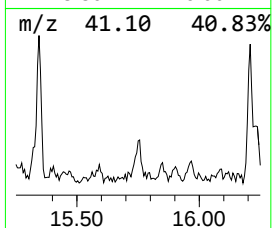
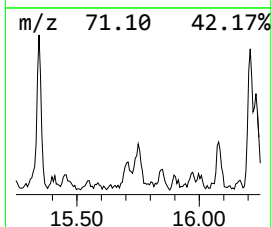
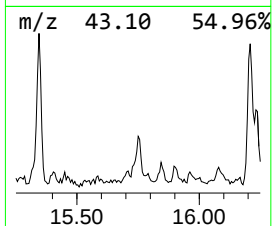
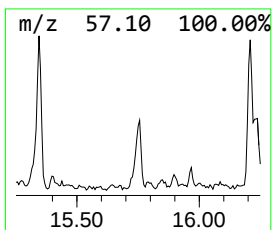
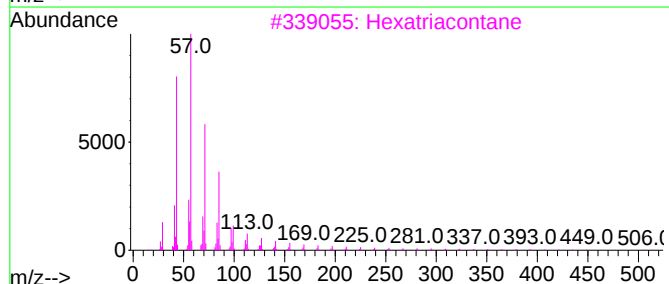
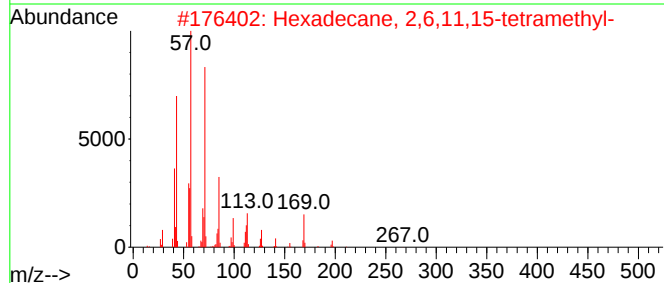
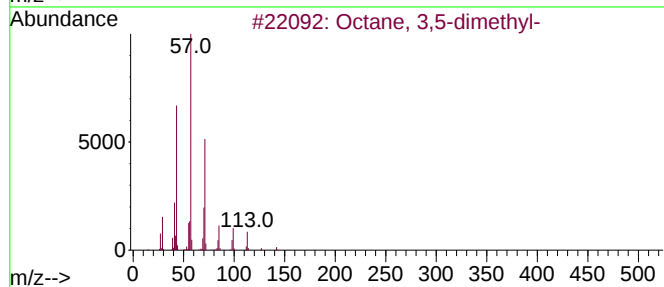
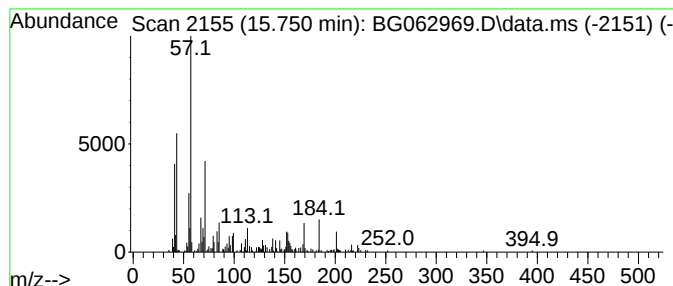
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Quant Title : SVOA CALIBRATION

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

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.750	8.80 ng/ul	264362	Acenaphthene-d10	14.487

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	60
2		Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	49
3		Hexatriacontane	507	C36H74	000630-06-8	43
4		9-Heptadecanone	254	C17H34O	000540-08-9	43
5		Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
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Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

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ClientSampleId :
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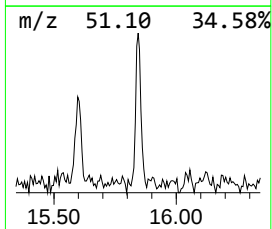
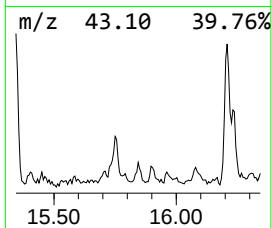
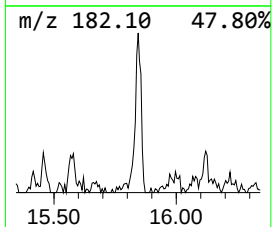
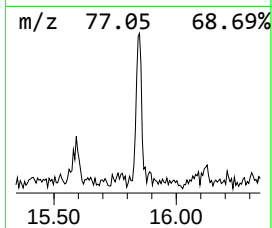
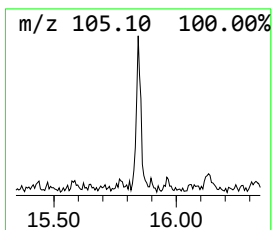
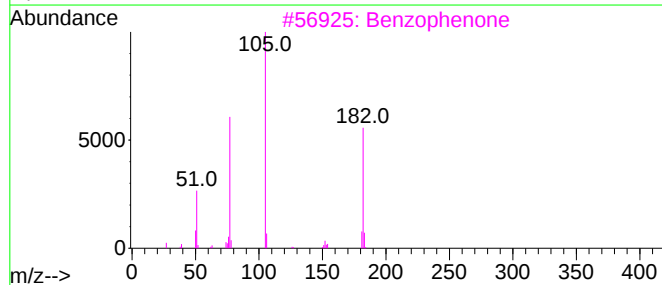
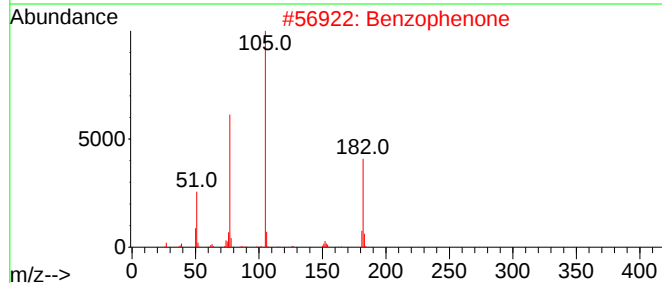
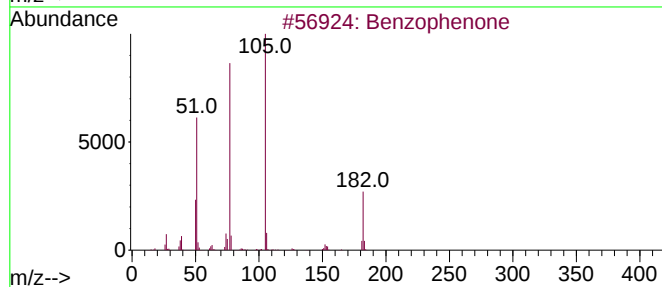
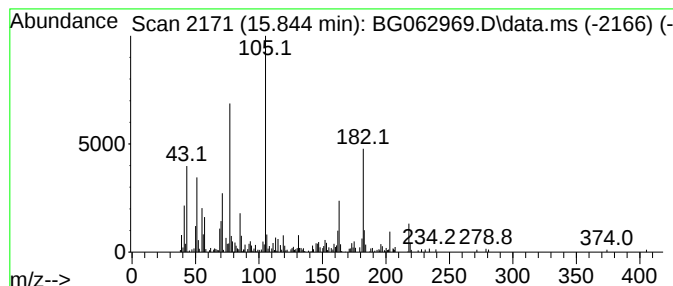
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 Benzophenone Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.844	7.56 ng/ul	227153	Acenaphthene-d10	14.487

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	78
2			Benzophenone	182	C13H10O	000119-61-9	60
3			Benzophenone	182	C13H10O	000119-61-9	53
4			Benzophenone	182	C13H10O	000119-61-9	53
5			Benzophenone	182	C13H10O	000119-61-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

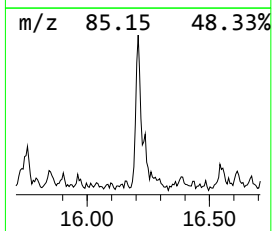
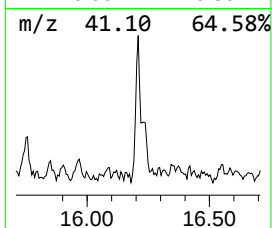
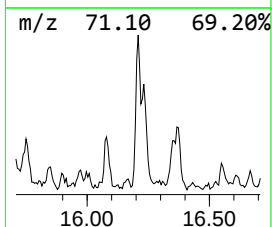
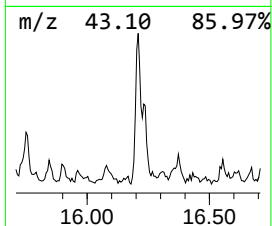
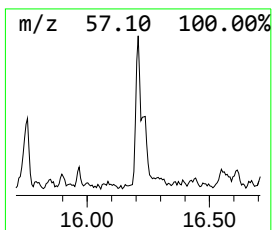
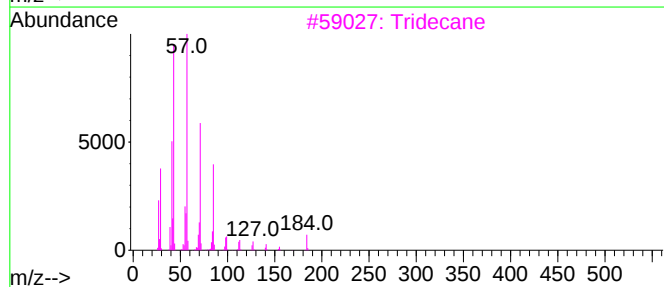
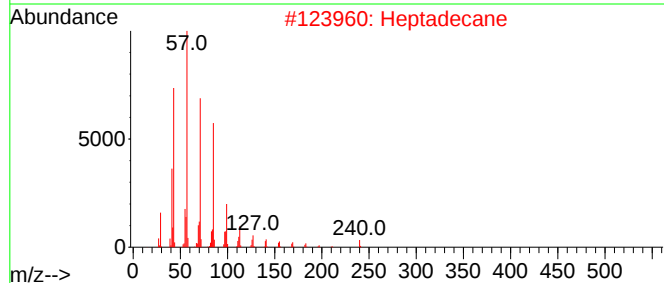
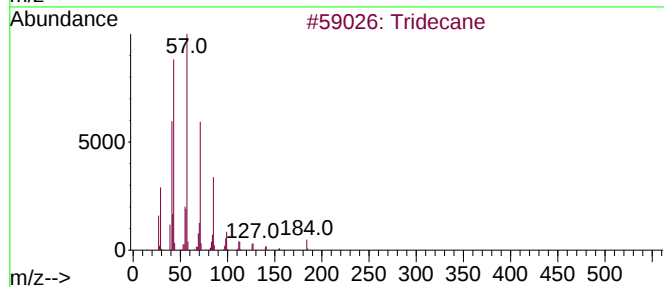
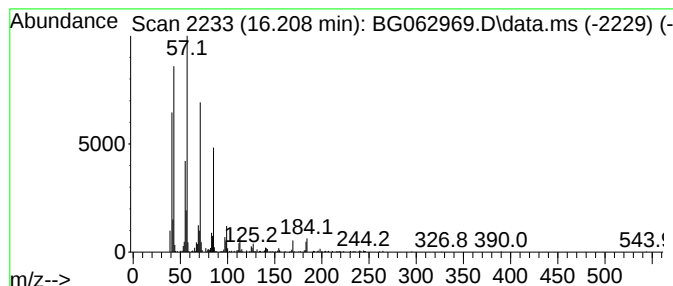
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BNA_G
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.208	13.42 ng/ul	524898	Phenanthrene-d10		17.242	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	95
2	Heptadecane		240	C17H36	000629-78-7	95
3	Tridecane		184	C13H28	000629-50-5	91
4	Tetradecane, 2,6,10-trimethyl-		240	C17H36	014905-56-7	90
5	Tridecane		184	C13H28	000629-50-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

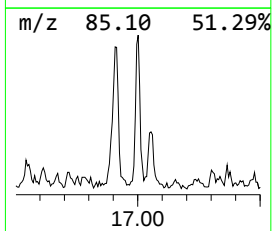
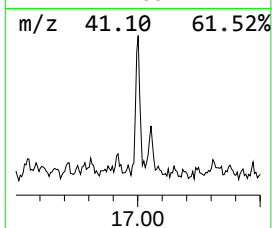
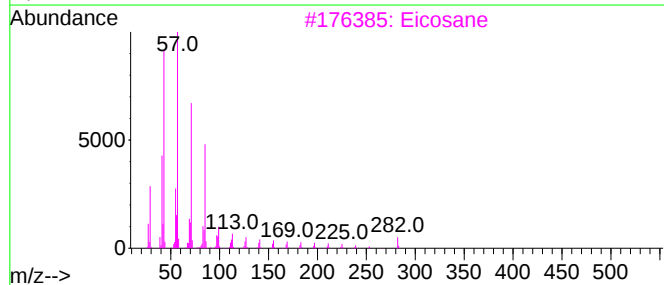
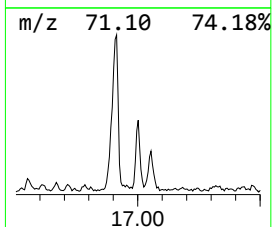
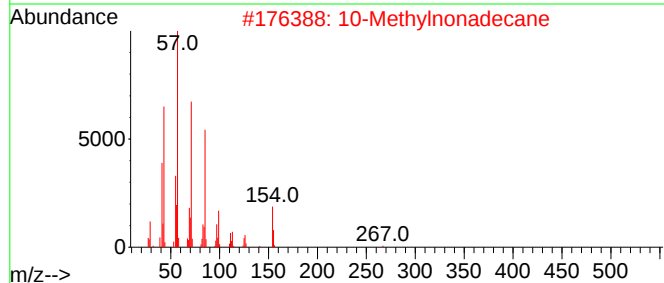
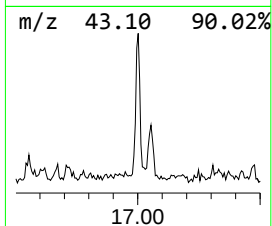
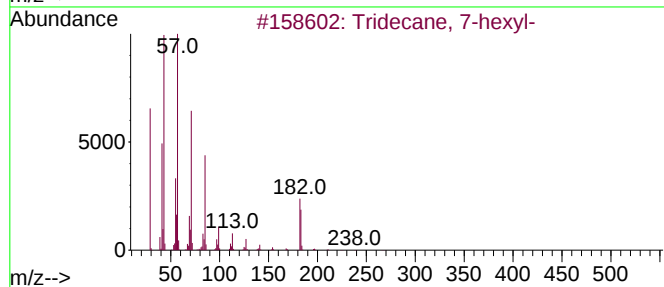
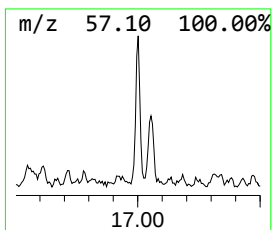
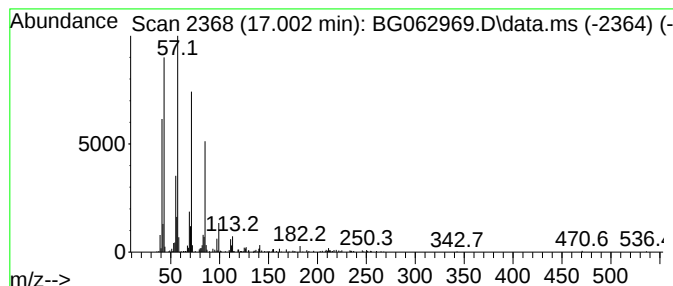
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.002	11.61 ng/ul	454126	Phenanthrene-d10		17.242	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane, 7-hexyl-		268	C19H40	007225-66-3	91
2	10-Methylnonadecane		282	C20H42	056862-62-5	91
3	Eicosane		282	C20H42	000112-95-8	91
4	Octadecane, 2-methyl-		268	C19H40	001560-88-9	90
5	Tetratetracontane		619	C44H90	007098-22-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

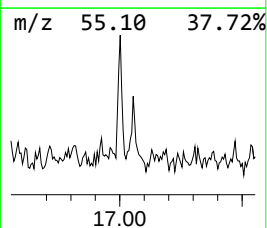
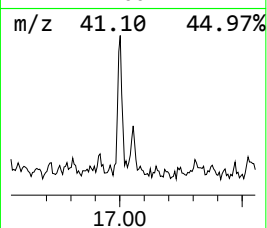
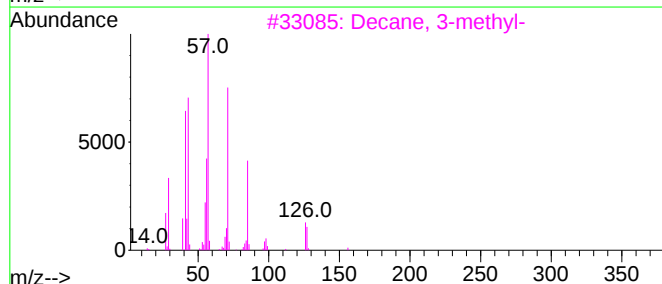
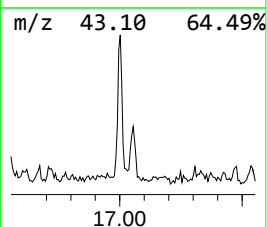
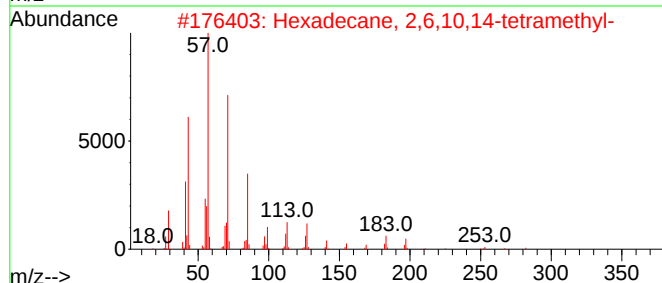
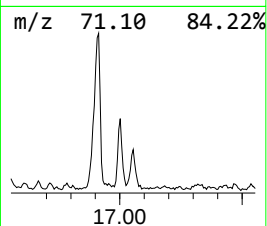
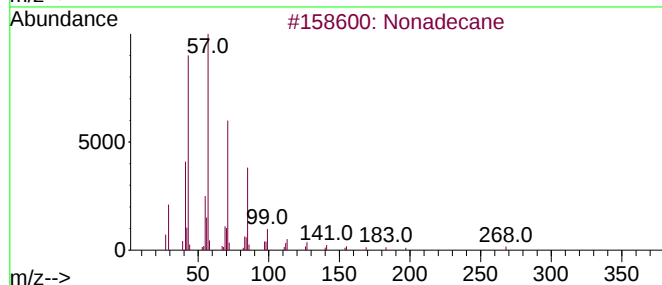
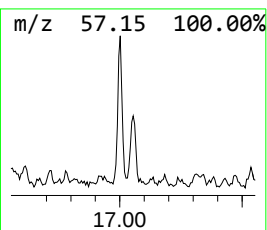
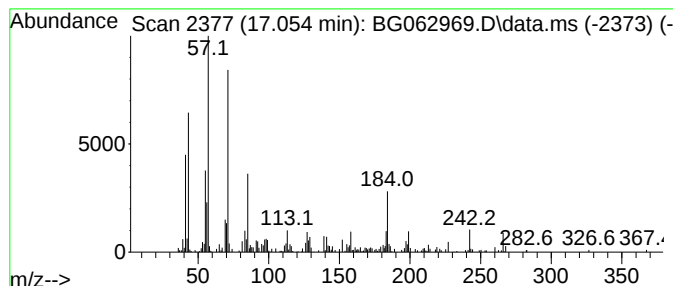
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.054	8.51 ng/ul	332833	Phenanthrene-d10	17.242

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	70
2		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	60
3		Decane, 3-methyl-	156	C11H24	013151-34-3	60
4		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	60
5		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

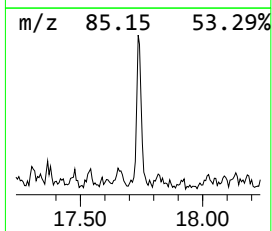
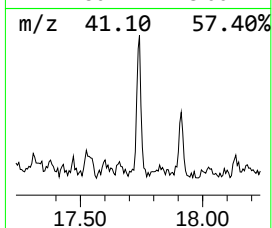
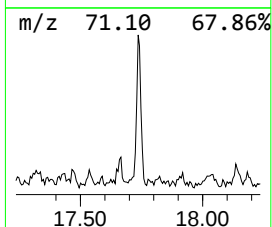
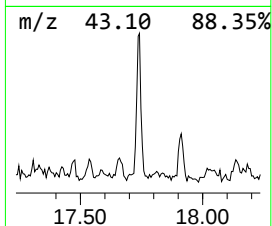
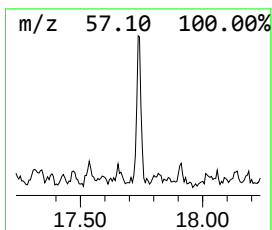
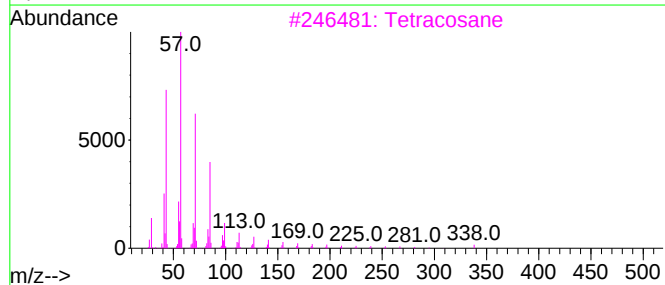
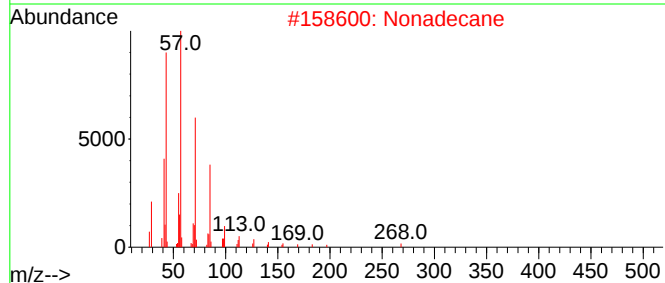
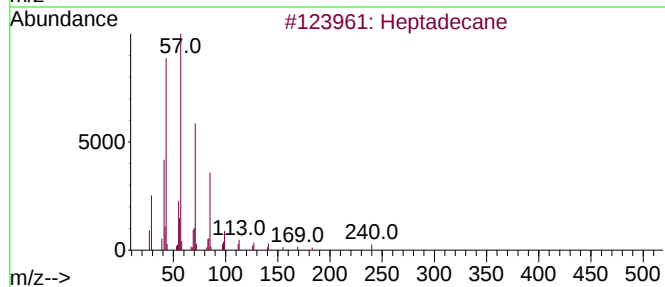
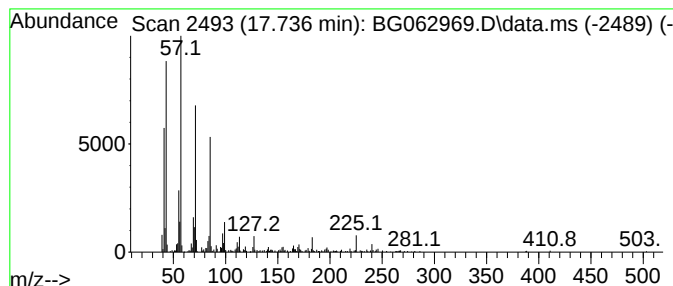
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.736	9.18 ng/ul	359072	Phenanthrene-d10	17.242

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	95
2		Nonadecane	268	C19H40	000629-92-5	83
3		Tetracosane	338	C24H50	000646-31-1	83
4		Heptadecane	240	C17H36	000629-78-7	83
5		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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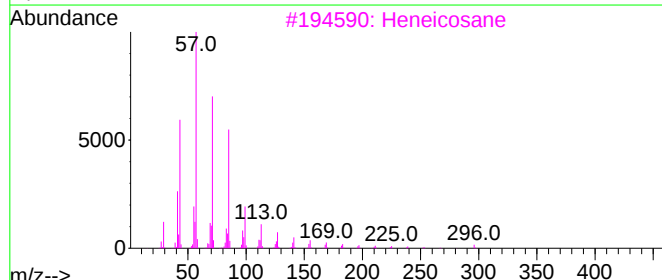
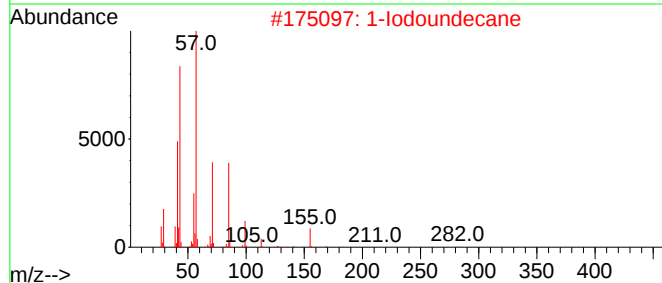
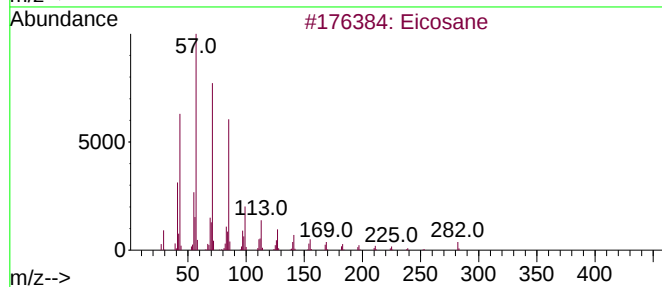
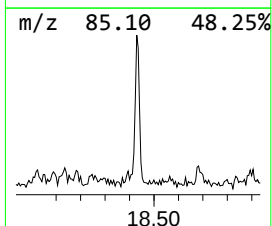
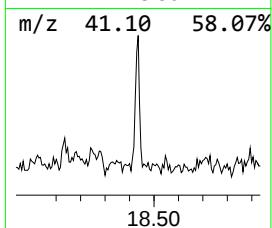
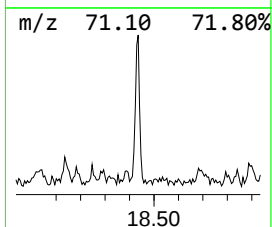
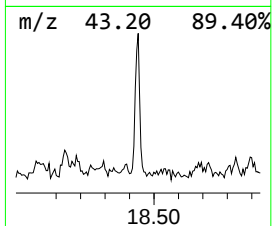
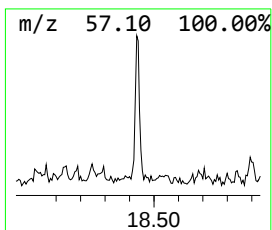
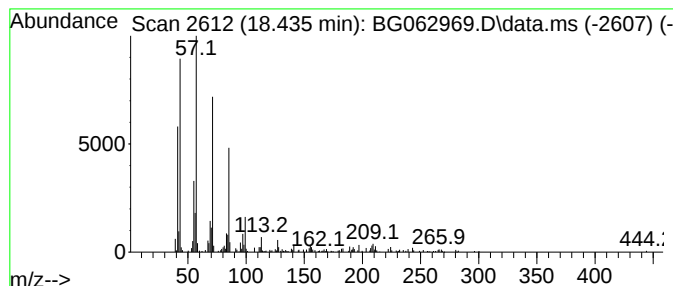
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Quant Title : SVOA CALIBRATION

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

Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.435	6.43 ng/ul	251393	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	91
2			1-Iodoundecane	282	C11H23I	004282-44-4	87
3			Heneicosane	296	C21H44	000629-94-7	83
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	83
5			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

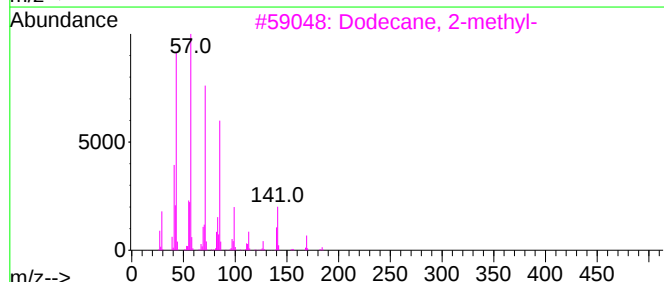
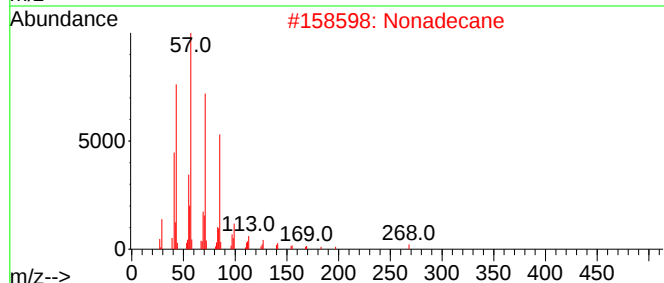
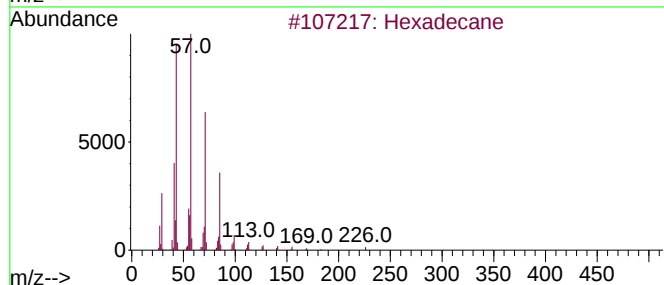
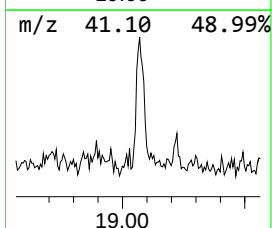
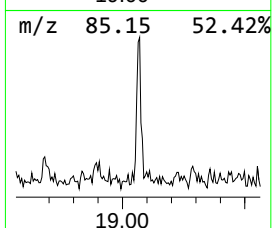
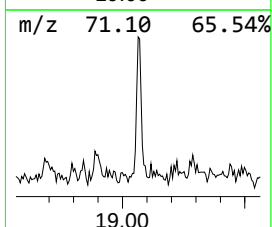
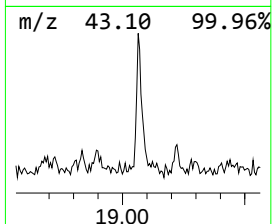
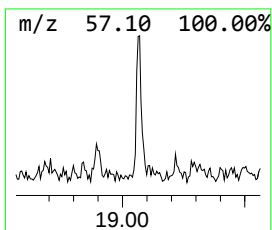
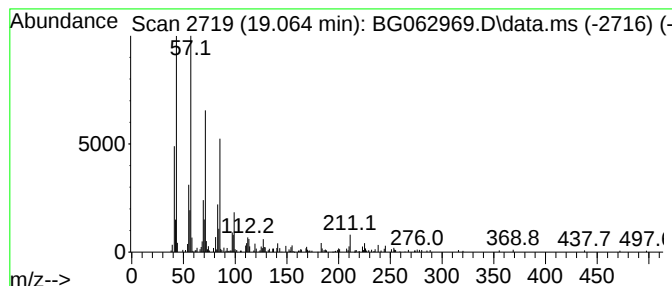
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.064	4.73 ng/ul	185076	Phenanthrene-d10	17.242

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	96
2			Nonadecane	268	C19H40	000629-92-5	87
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	87
4			Eicosane	282	C20H42	000112-95-8	83
5			Heneicosane	296	C21H44	000629-94-7	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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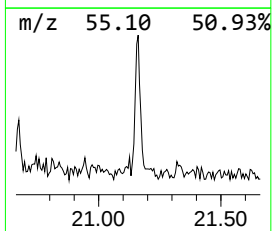
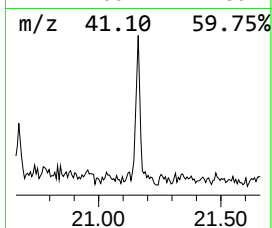
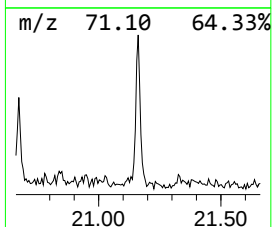
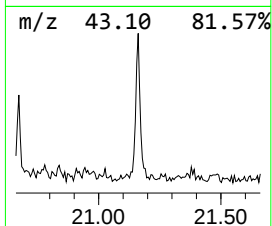
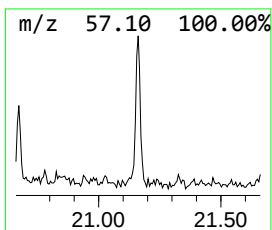
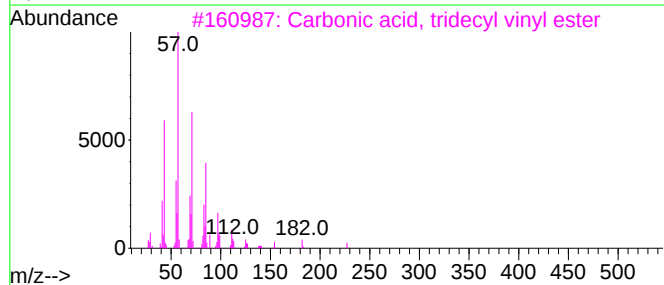
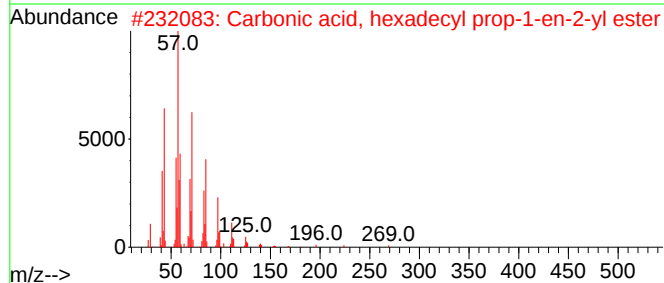
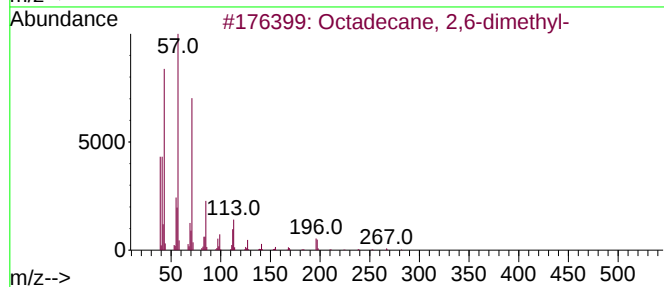
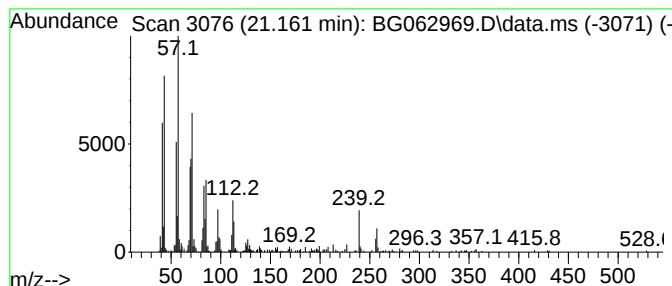
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Quant Title : SVOA CALIBRATION

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

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.161	8.37 ng/ul	331639	Chrysene-d12	21.485

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	51
2			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	49
3			Carbonic acid, tridecyl vinyl ester	270	C16H30O3	1000382-54-7	46
4			Octadecyl octyl ether	382	C26H54O	1000406-38-7	46
5			1-Tetradecene	196	C14H28	001120-36-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
Data File : BG062969.D
Acq On : 20 Sep 2024 8:39
Operator : JU/RC
Sample : P3898-20ME
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC59ME

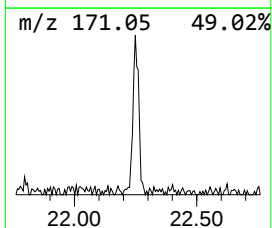
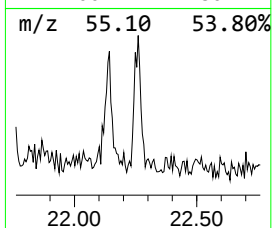
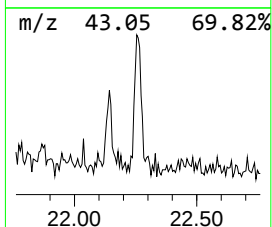
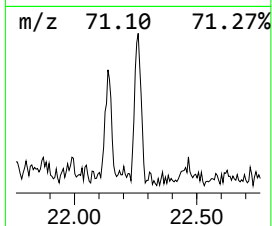
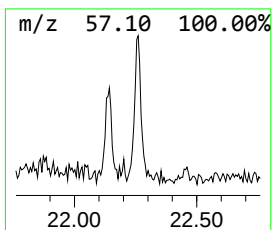
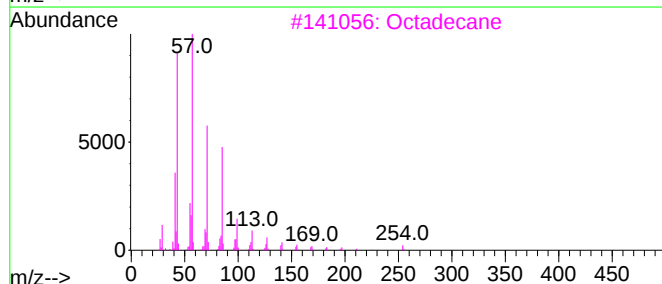
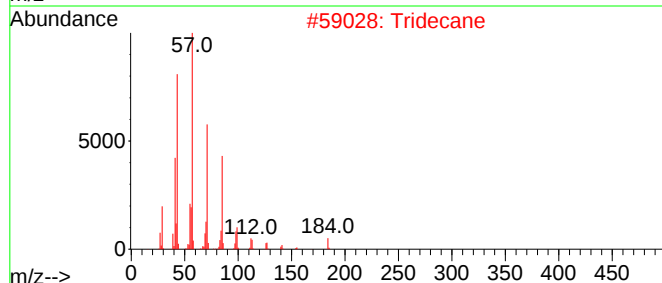
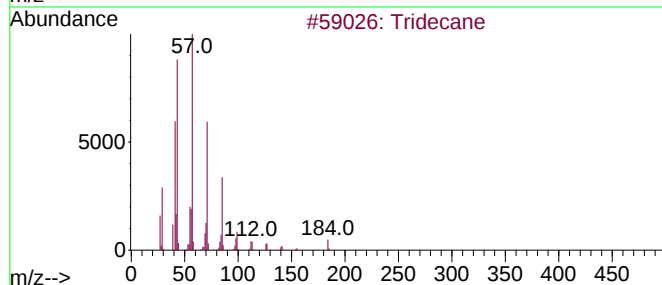
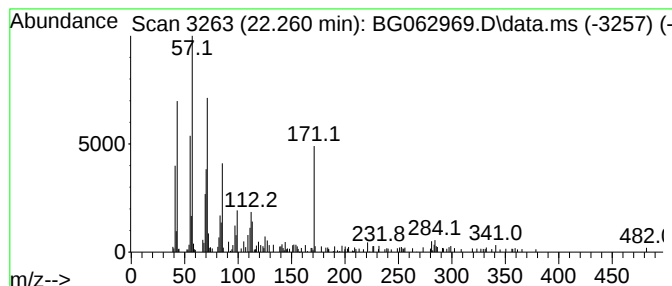
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Quant Title : SVOA CALIBRATION

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

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.260	4.70 ng/ul	185980	Chrysene-d12	21.485

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	50
2			Tridecane	184	C13H28	000629-50-5	50
3			Octadecane	254	C18H38	000593-45-3	46
4			Tridecane	184	C13H28	000629-50-5	46
5			Hexadecane, 7,9-dimethyl-	254	C18H38	021164-95-4	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091824\
 Data File : BG062969.D
 Acq On : 20 Sep 2024 8:39
 Operator : JU/RC
 Sample : P3898-20ME
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC59ME

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.885	19.7	ng/ul	528884	1	7.853	537867	20.0
unknown-01	6.355	6.3	ng/ul	169180	1	7.853	537867	20.0
(DEL) Alkane: S...	6.426	5.7	ng/ul	153280	1	7.853	537867	20.0
(DEL) Alkane: C...	6.508	6.8	ng/ul	183853	1	7.853	537867	20.0
(DEL) Alkane: S...	6.855	14.7	ng/ul	394369	1	7.853	537867	20.0
(DEL) Alkane: S...	6.913	9.8	ng/ul	263197	1	7.853	537867	20.0
Benzene, 1-ethy...	6.949	10.6	ng/ul	283686	1	7.853	537867	20.0
Benzene, 1,2,3-...	7.084	6.8	ng/ul	181416	1	7.853	537867	20.0
Benzene, 1-ethy...	7.248	7.3	ng/ul	197217	1	7.853	537867	20.0
2-Heptanone, 4,...	7.284	8.6	ng/ul	230636	1	7.853	537867	20.0
(DEL) Alkane: S...	7.489	65.4	ng/ul	1758430	1	7.853	537867	20.0
unknown-02	7.777	6.2	ng/ul	166897	1	7.853	537867	20.0
1-Hexanol, 2-et...	7.924	23.4	ng/ul	628128	1	7.853	537867	20.0
Mesitylene	7.971	9.2	ng/ul	247405	1	7.853	537867	20.0
Hexanal ethyl t...	8.077	16.8	ng/ul	450920	1	7.853	537867	20.0
(DEL) Alkane: C...	8.147	5.4	ng/ul	144398	1	7.853	537867	20.0
Cyclohexanone, ...	8.276	12.2	ng/ul	328962	1	7.853	537867	20.0
unknown-03	8.394	11.0	ng/ul	295068	1	7.853	537867	20.0
1,3,8-p-Menthat...	8.500	10.6	ng/ul	284869	1	7.853	537867	20.0
(DEL) Alkane: S...	8.623	10.0	ng/ul	268178	1	7.853	537867	20.0
Benzene, 1-meth...	8.658	8.5	ng/ul	229517	1	7.853	537867	20.0
Benzene, 2-ethy...	8.811	5.8	ng/ul	157265	1	7.853	537867	20.0
o-Cymene	8.858	9.3	ng/ul	250185	1	7.853	537867	20.0
(DEL) Alkane: S...	9.087	56.2	ng/ul	1512280	1	7.853	537867	20.0
unknown-04	9.181	7.2	ng/ul	192681	1	7.853	537867	20.0
unknown-05	9.205	9.0	ng/ul	241044	1	7.853	537867	20.0
(DEL) Alkane: S...	9.933	3.2	ng/ul	282768	2	10.644	1783370	20.0
(DEL) Alkane: S...	10.074	2.1	ng/ul	187841	2	10.644	1783370	20.0
2,4-Dipropyl-5-...	11.026	6.3	ng/ul	559565	2	10.644	1783370	20.0
(DEL) Alkane: S...	11.150	3.9	ng/ul	347145	2	10.644	1783370	20.0
(DEL) Alkane: S...	12.072	6.1	ng/ul	540337	2	10.644	1783370	20.0
(DEL) Alkane: S...	12.107	2.7	ng/ul	241765	2	10.644	1783370	20.0
(DEL) Alkane: S...	12.895	4.9	ng/ul	145934	3	14.487	600749	20.0
(DEL) Alkane: S...	13.318	19.3	ng/ul	580623	3	14.487	600749	20.0
Pyridine-2,4-di...	13.535	6.9	ng/ul	207502	3	14.487	600749	20.0
Naphthalene, 1,...	13.670	9.6	ng/ul	288788	3	14.487	600749	20.0
Butanoic acid, ...	13.752	6.5	ng/ul	194762	3	14.487	600749	20.0
Naphthalene, 1,...	13.811	5.4	ng/ul	161762	3	14.487	600749	20.0
(DEL) Alkane: S...	13.976	9.1	ng/ul	274372	3	14.487	600749	20.0
unknown-06	14.123	7.0	ng/ul	210827	3	14.487	600749	20.0
Sulfurous acid,...	14.393	31.3	ng/ul	939274	3	14.487	600749	20.0
unknown-07	14.563	7.6	ng/ul	227753	3	14.487	600749	20.0
Cyclopropane, 1...	14.628	16.6	ng/ul	498979	3	14.487	600749	20.0
Naphthalene, 2,...	14.886	7.4	ng/ul	222657	3	14.487	600749	20.0
(DEL) Alkane: C...	15.022	5.2	ng/ul	155646	3	14.487	600749	20.0
4,6,8-Trimethyl...	15.251	18.2	ng/ul	545521	3	14.487	600749	20.0
Tetradecane, 1-...	15.345	19.4	ng/ul	583395	3	14.487	600749	20.0
(DEL) Alkane: S...	15.750	8.8	ng/ul	264362	3	14.487	600749	20.0
Benzophenone	15.844	7.6	ng/ul	227153	3	14.487	600749	20.0
(DEL) Alkane: S...	16.208	13.4	ng/ul	524898	4	17.242	782395	20.0
(DEL) Alkane: S...	17.002	11.6	ng/ul	454126	4	17.242	782395	20.0
(DEL) Alkane: S...	17.054	8.5	ng/ul	332833	4	17.242	782395	20.0

(DEL) Alkane: S...	17.736	9.2	ng/ul	359072	4	17.242	782395	20.0
(DEL) Alkane: S...	18.435	6.4	ng/ul	251393	4	17.242	782395	20.0
(DEL) Alkane: S...	19.064	4.7	ng/ul	185076	4	17.242	782395	20.0
(DEL) Alkane: S...	21.161	8.4	ng/ul	331639	5	21.485	792120	20.0
(DEL) Alkane: S...	22.260	4.7	ng/ul	185980	5	21.485	792120	20.0

Instrument :

BNA_G

ClientSampleId :

DDC59ME