

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
 Data File : BG063109.D
 Acq On : 28 Sep 2024 15:45
 Operator : RC/JU
 Sample : P3910-12DL2 20X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DDC32DL2

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-BG092524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BG063109.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.739	108	111	115	rBV3	8642	11412	0.29%	0.059%
2	4.063	162	166	172	rVB6	8601	13187	0.34%	0.069%
3	4.967	315	320	325	rBV2	33671	46185	1.18%	0.240%
4	5.878	470	475	477	rBV5	18279	24255	0.62%	0.126%
5	6.160	518	523	529	rBV	84931	133409	3.42%	0.695%
6	6.236	532	536	540	rBV4	9026	16141	0.41%	0.084%
7	6.342	549	554	556	rBV3	10274	13121	0.34%	0.068%
8	6.413	564	566	571	rVB3	9062	11909	0.30%	0.062%
9	6.848	635	640	644	rVB6	11690	18850	0.48%	0.098%
10	6.906	646	650	653	rBV3	13955	17755	0.45%	0.092%
11	6.942	653	656	661	rVB2	12002	18421	0.47%	0.096%
12	7.000	661	666	667	rBV2	24076	29032	0.74%	0.151%
13	7.077	676	679	684	rVB2	13498	19844	0.51%	0.103%
14	7.177	691	696	702	rBV6	15169	30924	0.79%	0.161%
15	7.276	709	713	719	rVV4	21102	38886	1.00%	0.202%
16	7.382	727	731	737	rVB3	13751	26271	0.67%	0.137%
17	7.482	742	748	756	rVV3	52087	120731	3.09%	0.629%
18	7.846	804	810	817	rBV	260636	437363	11.20%	2.277%
19	7.911	817	821	827	rVB3	27720	49503	1.27%	0.258%
20	8.140	859	860	866	rVB4	10728	12018	0.31%	0.063%
21	8.275	877	883	887	rBV2	111079	181921	4.66%	0.947%
22	8.399	898	904	908	rVB7	17749	38136	0.98%	0.199%
23	8.469	908	916	926	rBV9	37310	140197	3.59%	0.730%
24	8.557	926	931	937	rVV6	19880	42172	1.08%	0.220%
25	8.616	937	941	944	rVV3	16198	26341	0.67%	0.137%
26	8.645	944	946	952	rVB2	17369	28562	0.73%	0.149%
27	8.798	967	972	975	rBV4	16638	23459	0.60%	0.122%
28	8.839	975	979	988	rVB6	31650	70092	1.79%	0.365%
29	8.963	988	1000	1004	rBV7	31218	77812	1.99%	0.405%
30	9.004	1005	1007	1012	rVV4	13216	20904	0.54%	0.109%
31	9.074	1015	1019	1024	rVB3	55508	93095	2.38%	0.485%
32	9.433	1077	1080	1083	rVB5	9483	11416	0.29%	0.059%
33	9.468	1083	1086	1091	rBV5	8453	14679	0.38%	0.076%
34	9.697	1120	1125	1126	rBV3	14462	15098	0.39%	0.079%
35	9.779	1136	1139	1145	rVB6	9794	11790	0.30%	0.061%
36	9.997	1172	1176	1180	rBV6	10764	20315	0.52%	0.106%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
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37	10.267	1217	1222	1229	rVB4	26428	52098	1.33%	0.271%
38	10.332	1229	1233	1236	rBV4	21704	26749	0.68%	0.139%
39	10.508	1258	1263	1268	rBV7	14763	26864	0.69%	0.140%
40	10.631	1275	1284	1296	rBV2	711935	1354411	34.67%	7.051%
41	10.755	1299	1305	1311	rBV3	126368	231154	5.92%	1.203%
42	10.931	1329	1335	1340	rBV6	17592	32987	0.84%	0.172%
43	11.019	1344	1350	1356	rVB4	44920	84058	2.15%	0.438%
44	11.113	1362	1366	1369	rBV6	14915	21391	0.55%	0.111%
45	11.148	1370	1372	1378	rVB5	18996	23551	0.60%	0.123%
46	11.536	1432	1438	1439	rBV4	26069	45645	1.17%	0.238%
47	11.665	1457	1460	1468	rBV6	14338	27644	0.71%	0.144%
48	11.742	1468	1473	1480	rVB3	31645	51914	1.33%	0.270%
49	11.883	1492	1497	1505	rBV7	20063	53221	1.36%	0.277%
50	12.059	1520	1527	1531	rBV5	50288	99214	2.54%	0.517%
51	12.312	1564	1570	1575	rVB6	28483	60506	1.55%	0.315%
52	12.482	1592	1599	1601	rVV5	22896	42386	1.09%	0.221%
53	12.523	1605	1606	1613	rVV6	15716	20336	0.52%	0.106%
54	12.888	1663	1668	1671	rBV6	7100	11799	0.30%	0.061%
55	13.052	1692	1696	1702	rVB2	43958	68768	1.76%	0.358%
56	13.311	1736	1740	1744	rBV3	30435	43396	1.11%	0.226%
57	13.534	1772	1778	1783	rVB10	17918	37565	0.96%	0.196%
58	13.657	1793	1799	1804	rVB8	22435	48773	1.25%	0.254%
59	13.810	1820	1825	1831	rVV7	26523	67212	1.72%	0.350%
60	13.863	1831	1834	1835	rVV3	21069	25629	0.66%	0.133%
61	13.886	1835	1838	1844	rVB	66846	97123	2.49%	0.506%
62	13.969	1849	1852	1856	rVB6	17519	21851	0.56%	0.114%
63	14.027	1856	1862	1868	rBV9	25962	51026	1.31%	0.266%
64	14.133	1873	1880	1883	rBV8	31077	64665	1.66%	0.337%
65	14.180	1883	1888	1892	rVV	62673	81382	2.08%	0.424%
66	14.386	1917	1923	1927	rBV3	51751	76967	1.97%	0.401%
67	14.427	1928	1930	1933	rVV4	18794	15367	0.39%	0.080%
68	14.486	1933	1940	1947	rVB	787808	1258712	32.22%	6.553%
69	14.627	1960	1964	1970	rVB2	59401	89545	2.29%	0.466%
70	14.703	1972	1977	1980	rBV7	20140	41257	1.06%	0.215%
71	14.826	1996	1998	2003	rVB6	12320	13687	0.35%	0.071%
72	14.956	2018	2020	2025	rVB5	12955	16492	0.42%	0.086%
73	15.009	2025	2029	2030	rBV4	15028	14384	0.37%	0.075%
74	15.114	2042	2047	2056	rVB2	219253	344033	8.81%	1.791%
75	15.244	2066	2069	2074	rVV3	46582	66338	1.70%	0.345%
76	15.314	2078	2081	2082	rBV3	12445	14204	0.36%	0.074%

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77	15.343	2082	2086	2089	rVB2	39022	50834	1.30%	0.265%
78	15.479	2105	2109	2113	rVB	96783	128443	3.29%	0.669%
79	15.843	2167	2171	2177	rVB3	56597	95561	2.45%	0.498%
80	16.043	2201	2205	2209	rBV7	10695	17294	0.44%	0.090%
81	16.113	2214	2217	2223	rVB6	25146	44055	1.13%	0.229%
82	16.207	2229	2233	2236	rBV2	55651	82891	2.12%	0.432%
83	16.542	2287	2290	2294	rVB6	22662	34146	0.87%	0.178%
84	16.889	2343	2349	2360	rBV	2554832	3906500	100.00%	20.338%
85	16.994	2365	2367	2371	rVB5	45382	51323	1.31%	0.267%
86	17.235	2402	2408	2417	rBV	1279497	1876441	48.03%	9.769%
87	17.335	2420	2425	2430	rVB2	116877	178388	4.57%	0.929%
88	17.629	2471	2475	2476	rBV4	9973	11762	0.30%	0.061%
89	17.735	2490	2493	2498	rVB4	51012	75225	1.93%	0.392%
90	17.811	2502	2506	2510	rBV3	28235	43494	1.11%	0.226%
91	18.428	2608	2611	2615	rBV3	41825	50802	1.30%	0.264%
92	18.516	2621	2626	2627	rBV5	12390	18929	0.48%	0.099%
93	19.063	2716	2719	2722	rBV4	21713	31762	0.81%	0.165%
94	19.633	2811	2816	2822	rVV	176495	270768	6.93%	1.410%
95	20.179	2906	2909	2910	rBV2	23100	24461	0.63%	0.127%
96	21.160	3073	3076	3082	rVB6	49231	72884	1.87%	0.379%
97	21.489	3125	3132	3140	rBV	1510179	2389890	61.18%	12.442%
98	22.905	3371	3373	3381	rVB9	33166	55258	1.41%	0.288%
99	24.315	3608	3613	3621	rVB2	88161	213128	5.46%	1.110%
100	24.527	3640	3649	3669	rVB	1001978	2754265	70.50%	14.339%

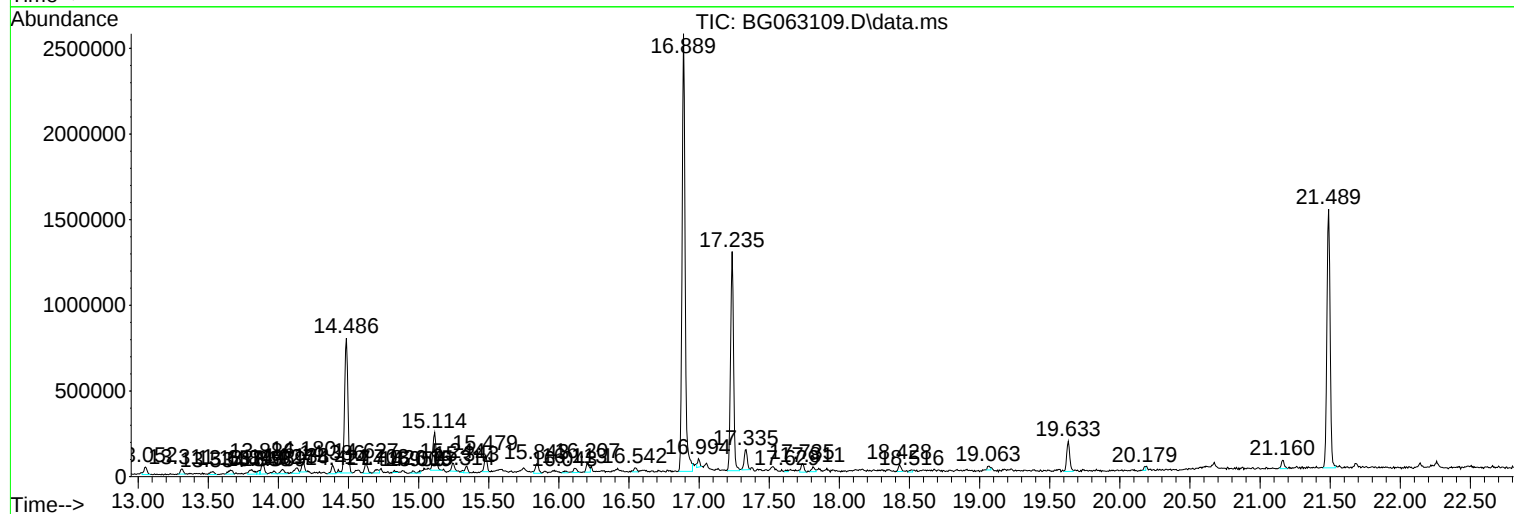
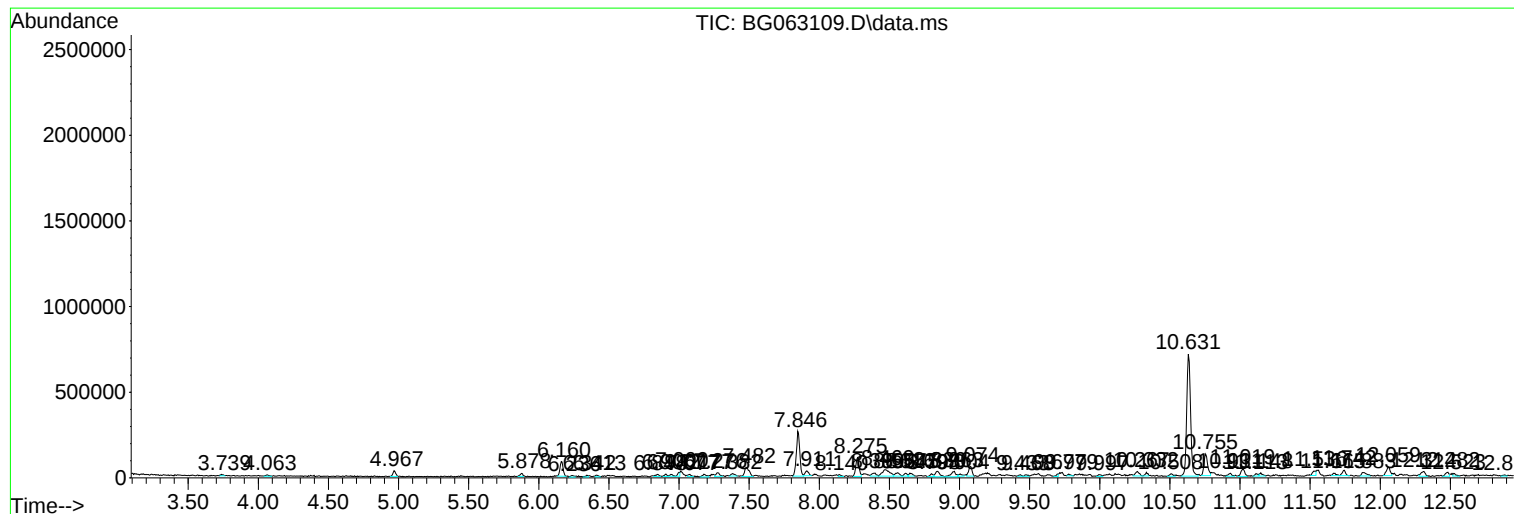
Sum of corrected areas: 19208014

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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



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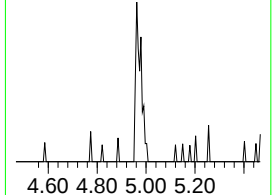
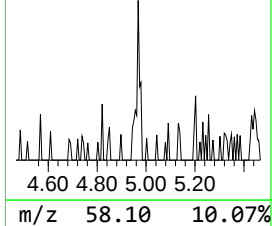
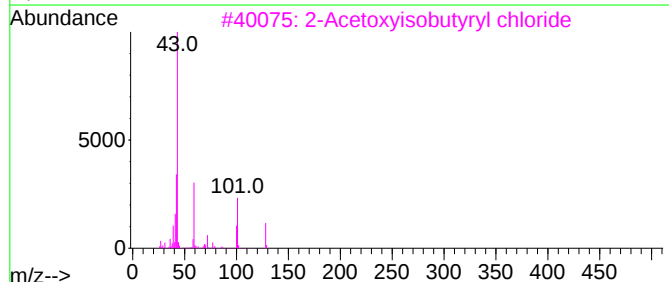
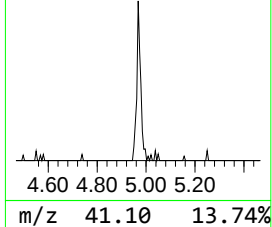
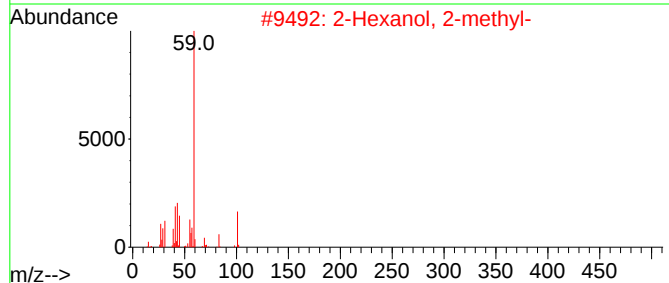
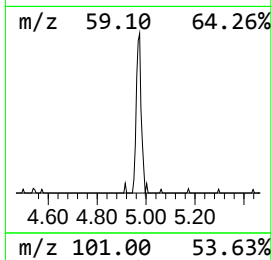
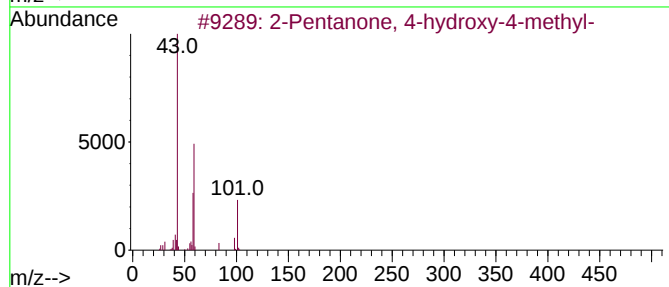
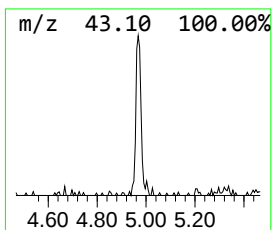
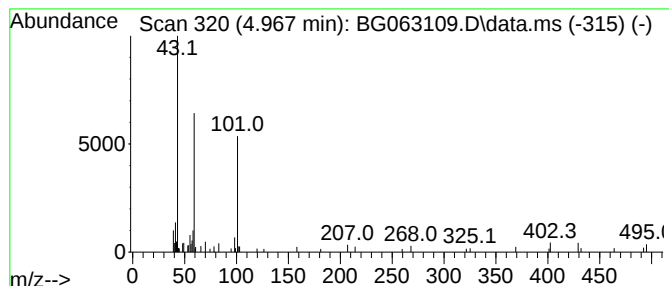
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.967	2.11 ng/ul	46185	1,4-Dichlorobenzene-d4	7.846

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53		
2	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	42		
3	2-Acetoxyisobutyryl chloride	164	C6H9ClO3	040635-66-3	42		
4	(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	37		
5	2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	36		



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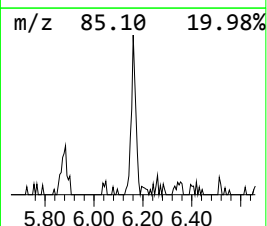
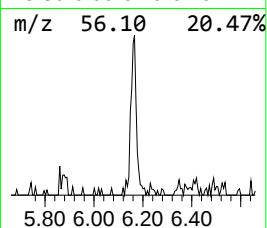
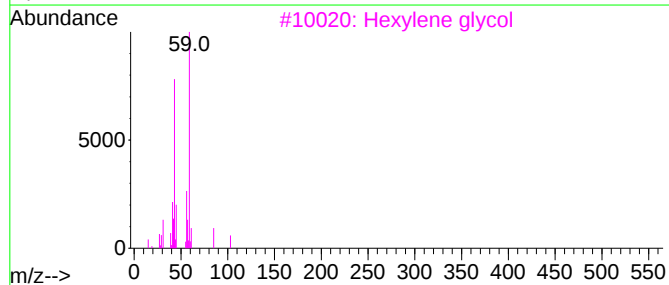
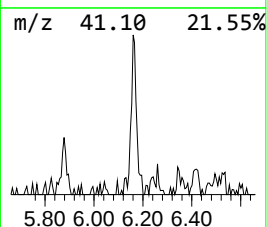
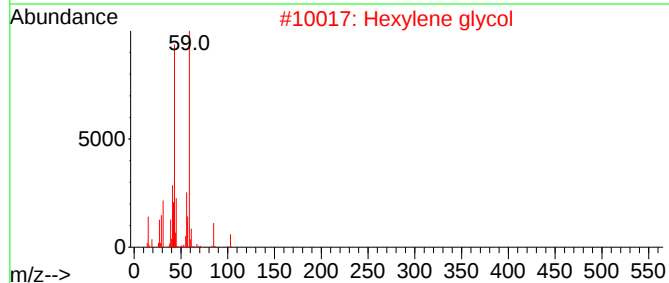
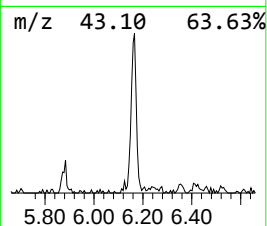
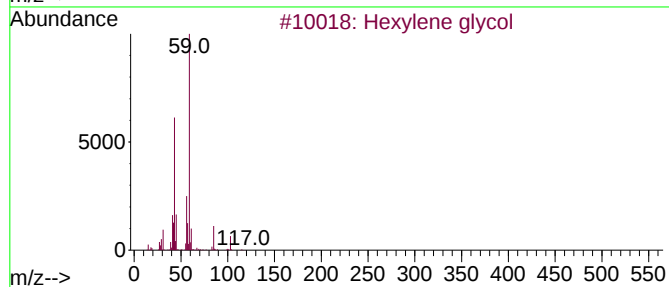
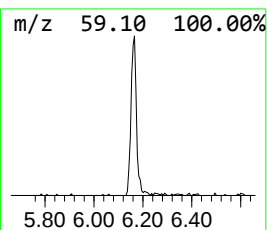
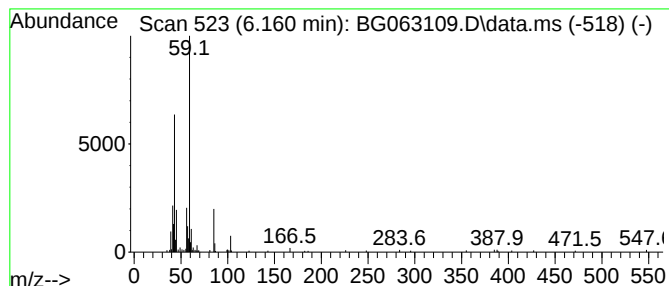
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Hexylene glycol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.160	6.10 ng/ul	133409	1,4-Dichlorobenzene-d4	7.846

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexylene glycol	118	C6H14O2	000107-41-5	72
2			Hexylene glycol	118	C6H14O2	000107-41-5	64
3			Hexylene glycol	118	C6H14O2	000107-41-5	64
4			Hexylene glycol	118	C6H14O2	000107-41-5	50
5			2-Propanol, 1-methoxy-2-methyl-	104	C5H12O2	003587-64-2	47



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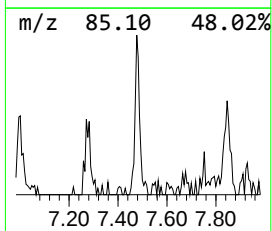
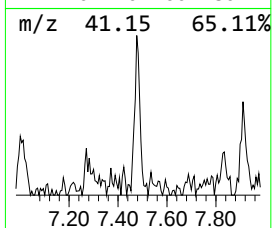
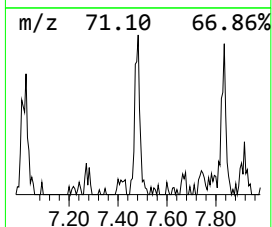
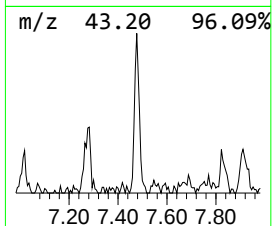
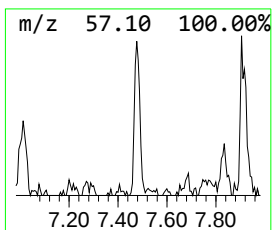
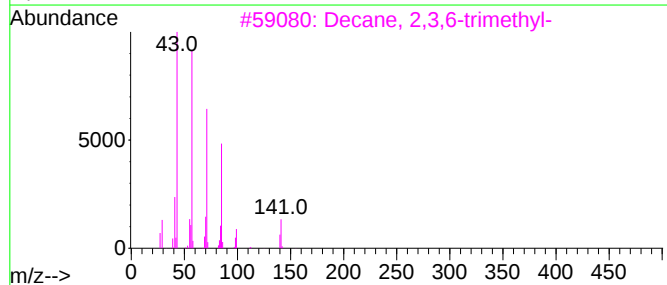
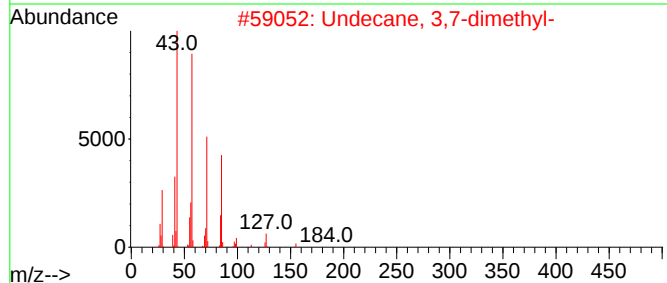
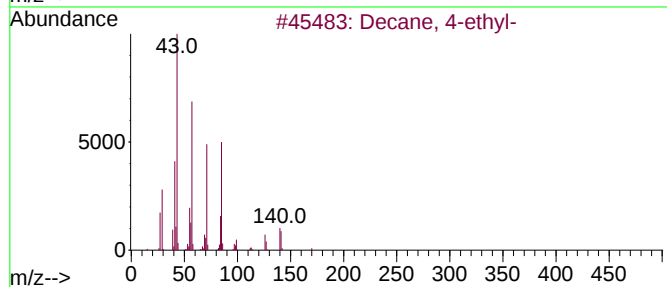
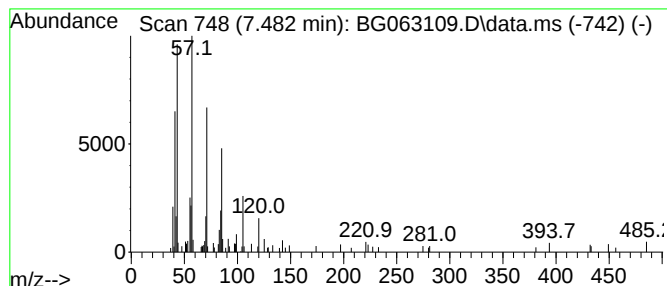
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.482	5.52 ng/ul	120731	1,4-Dichlorobenzene-d4			7.846
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 4-ethyl-		170	C12H26	001636-44-8	59
2	Undecane, 3,7-dimethyl-		184	C13H28	017301-29-0	59
3	Decane, 2,3,6-trimethyl-		184	C13H28	062238-12-4	50
4	Decane, 2,4-dimethyl-		170	C12H26	002801-84-5	50
5	4-Heptanone, 2-methyl-		128	C8H16O	000626-33-5	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063109.D
Acq On : 28 Sep 2024 15:45
Operator : RC/JU
Sample : P3910-12DL2 20X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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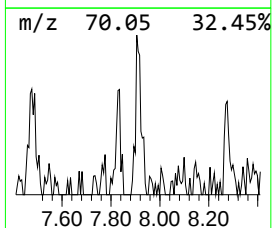
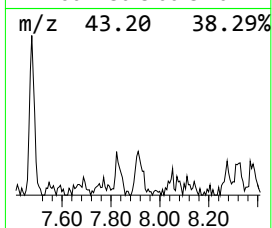
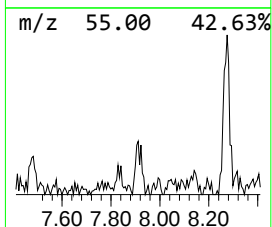
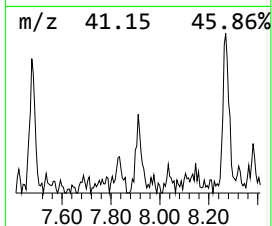
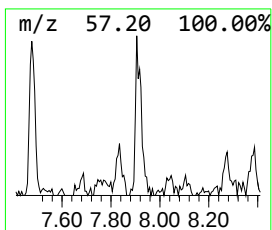
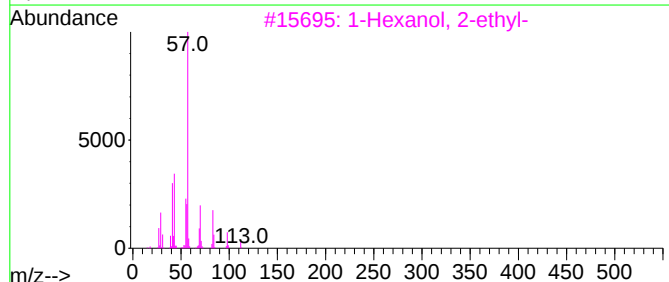
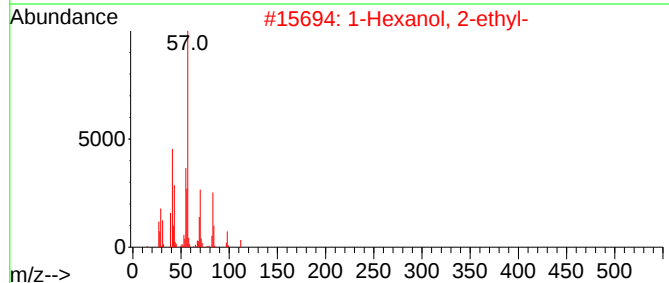
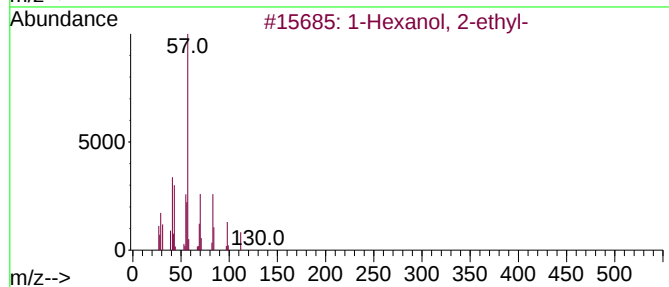
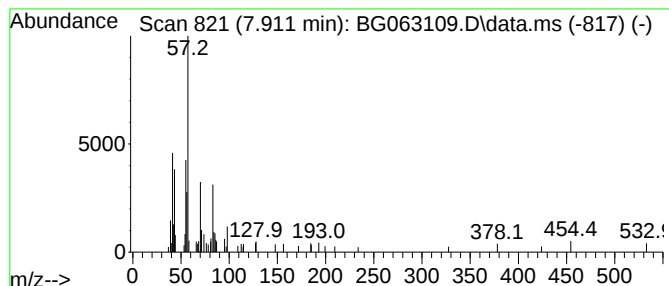
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.911	2.26 ng/ul	49503	1,4-Dichlorobenzene-d4	7.846

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	50
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	50
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	50
4	2-Propyl-1-pentanol			130	C8H18O	058175-57-8	43
5	1-Decene, 2,4-dimethyl-			168	C12H24	055170-80-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063109.D
Acq On : 28 Sep 2024 15:45
Operator : RC/JU
Sample : P3910-12DL2 20X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DDC32DL2

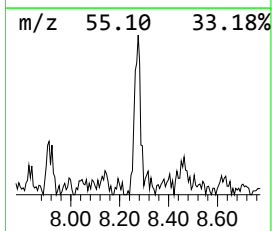
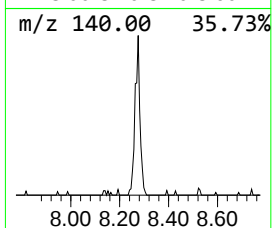
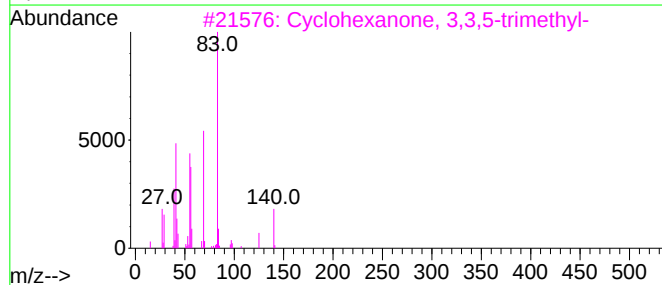
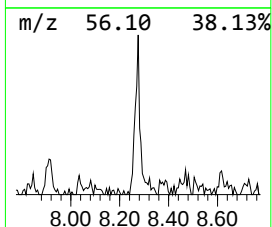
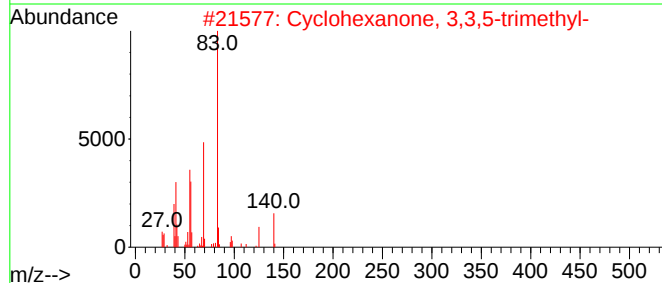
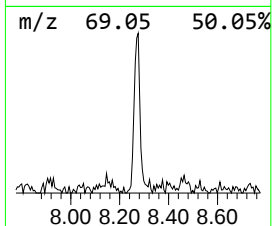
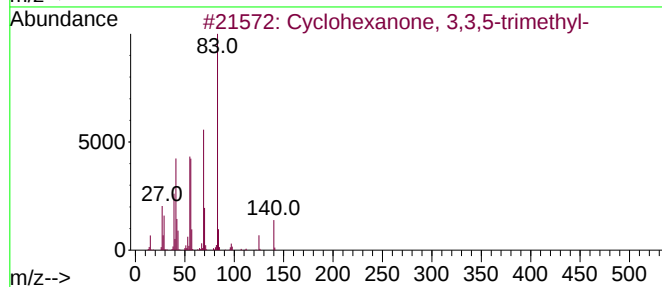
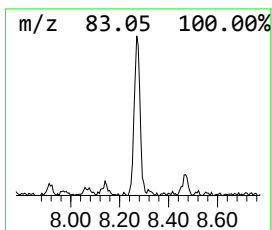
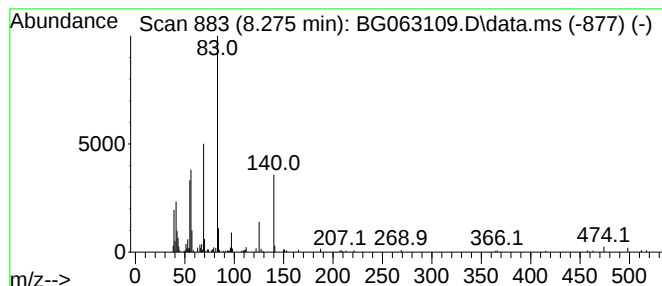
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclohexanone, 3,3,5-trimet... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.275	8.32 ng/ul	181921	1,4-Dichlorobenzene-d4	7.846

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	90
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	90
5			1H-1,2,4-Triazole, 3-methyl-	83	C3H5N3	007170-01-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063109.D
Acq On : 28 Sep 2024 15:45
Operator : RC/JU
Sample : P3910-12DL2 20X
Misc :
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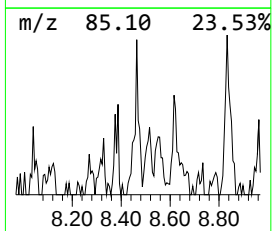
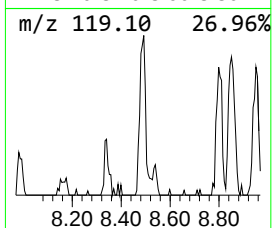
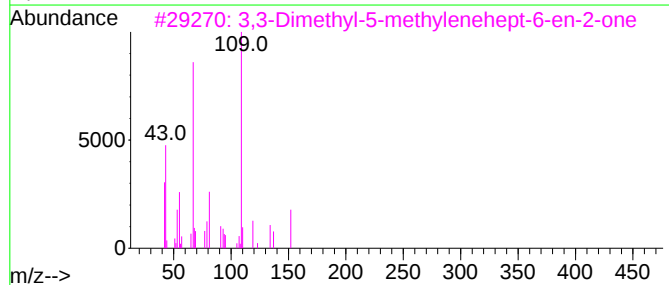
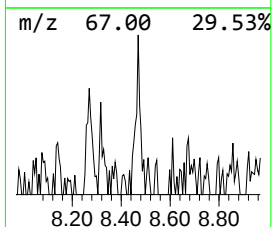
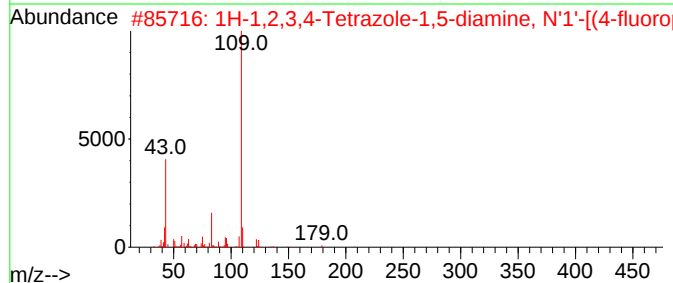
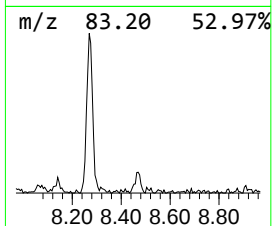
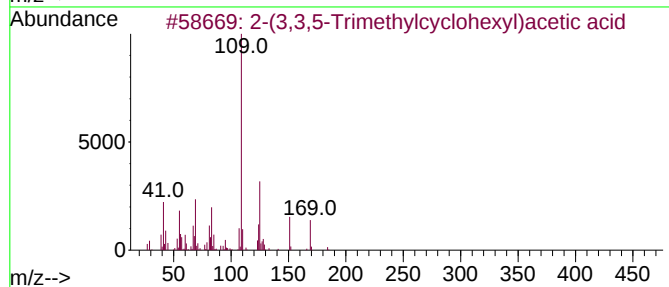
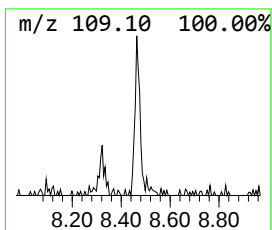
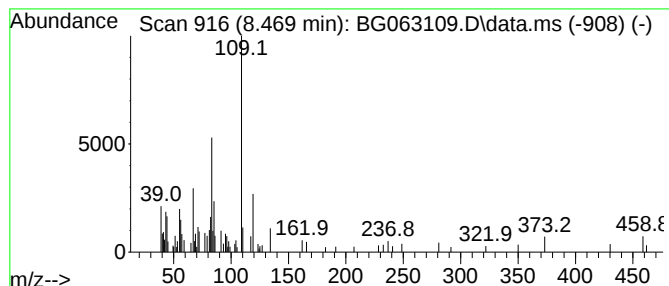
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-(3,3,5-Trimethylcyclohexy... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.469	6.41 ng/ul	140197	1,4-Dichlorobenzene-d4	7.846	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(3,3,5-Trimethylcyclohexyl)ace...	184	C11H20O2	003213-73-8	53
2	1H-1,2,3,4-Tetrazole-1,5-diamine...	208	C8H9FN6	1000351-63-3	38
3	3,3-Dimethyl-5-methylenehept-6-e...	152	C10H16O	1010424-30-5	35
4	(3-Fluorophenyl) methanol, tert...	182	C11H15FO	1000374-63-2	35
5	6,7-Dihydro-5H-benzo[1,2,5]oxadi...	261	C13H12FN3O2	1000300-92-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063109.D
Acq On : 28 Sep 2024 15:45
Operator : RC/JU
Sample : P3910-12DL2 20X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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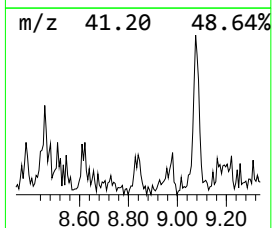
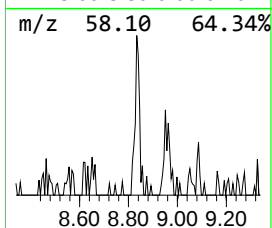
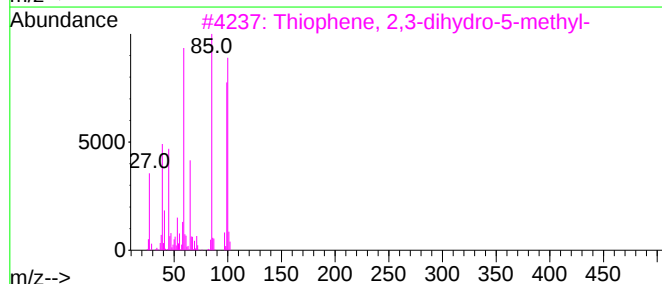
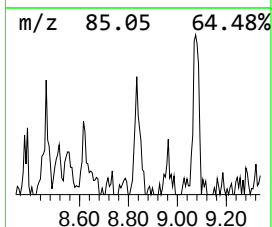
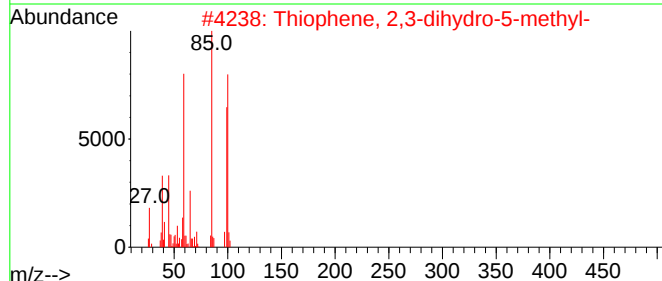
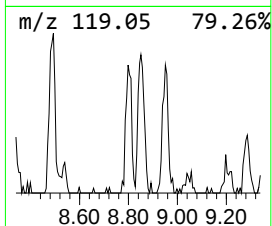
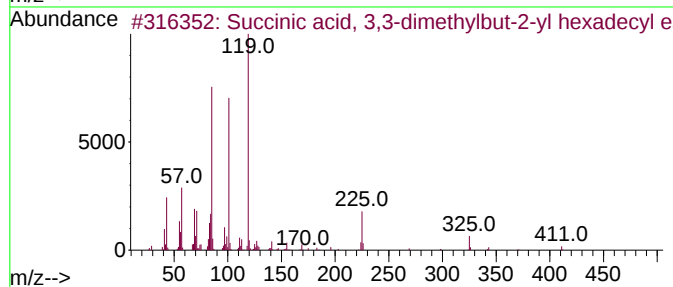
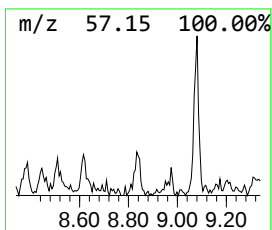
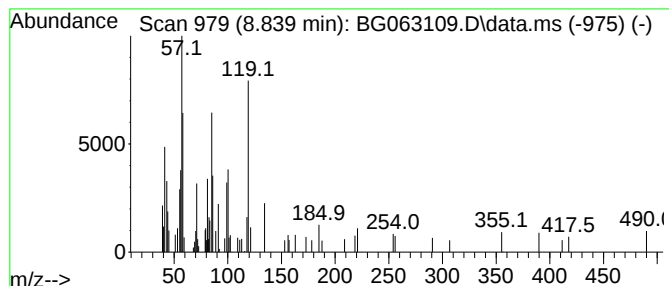
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.839	3.21 ng/ul	70092	1,4-Dichlorobenzene-d4	7.846

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Succinic acid, 3,3-dimethylbut-2...	426	C26H50O4	1000349-52-0	38
2			Thiophene, 2,3-dihydro-5-methyl-	100	C5H8S	004610-02-0	27
3			Thiophene, 2,3-dihydro-5-methyl-	100	C5H8S	004610-02-0	27
4			Thiophene, 2,3-dihydro-5-methyl-	100	C5H8S	004610-02-0	27
5			[6-N-(N,N-dimethylaminomethylene...	200	C10H20N2O2	1000370-32-1	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063109.D
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Operator : RC/JU
Sample : P3910-12DL2 20X
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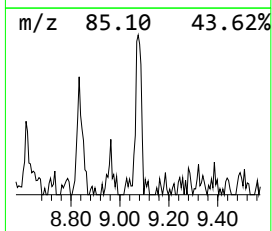
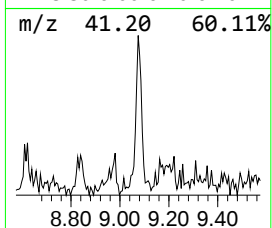
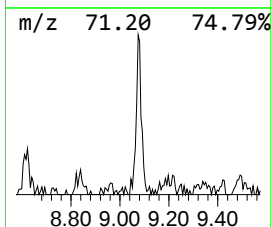
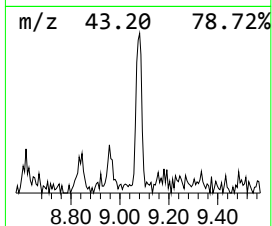
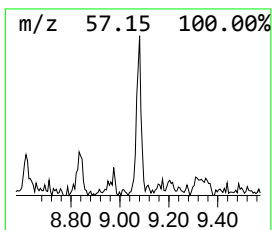
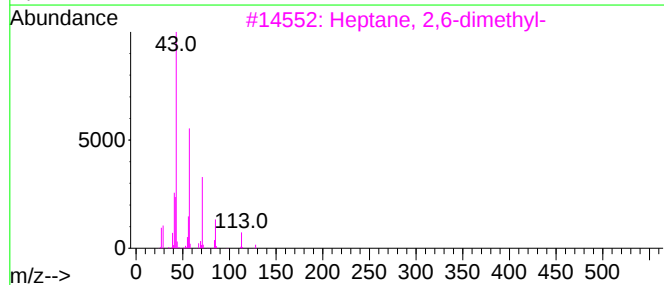
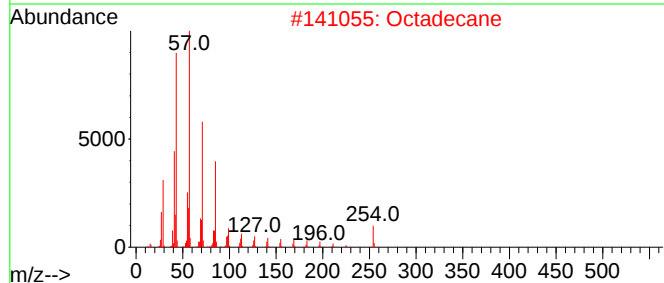
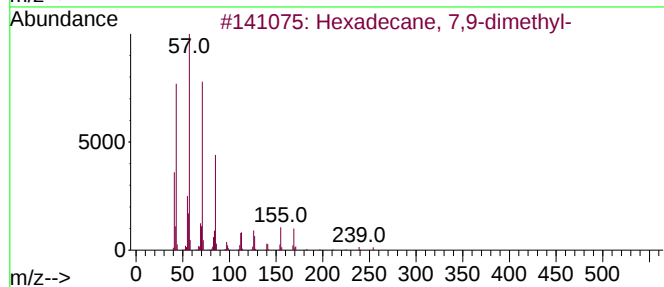
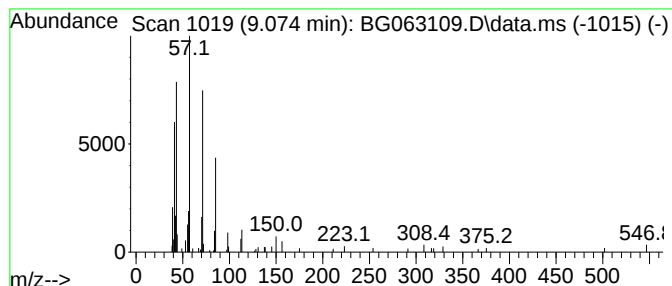
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.074	4.26 ng/ul	93095	1,4-Dichlorobenzene-d4			7.846
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane, 7,9-dimethyl-		254	C18H38	021164-95-4	87
2	Octadecane		254	C18H38	000593-45-3	78
3	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	76
4	Undecane		156	C11H24	001120-21-4	74
5	Octane, 2-methyl-		128	C9H20	003221-61-2	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092724\
Data File : BG063109.D
Acq On : 28 Sep 2024 15:45
Operator : RC/JU
Sample : P3910-12DL2 20X
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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
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Hexylene glycol	6.160	6.1	ng/ul	133409	1	7.846	437363	20.0
(DEL) Alkane: S...	7.482	5.5	ng/ul	120731	1	7.846	437363	20.0
1-Hexanol, 2-et...	7.911	2.3	ng/ul	49503	1	7.846	437363	20.0
Cyclohexanone, ...	8.275	8.3	ng/ul	181921	1	7.846	437363	20.0
2-(3,3,5-Trimet...	8.469	6.4	ng/ul	140197	1	7.846	437363	20.0
unknown-01	8.839	3.2	ng/ul	70092	1	7.846	437363	20.0
(DEL) Alkane: S...	9.074	4.3	ng/ul	93095	1	7.846	437363	20.0