

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022109.D
 Acq On : 28 Sep 2024 17:48
 Operator : RC/JU
 Sample : P3910-10MEDL 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC29MEDL

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_P\Methods\SFAM-EPA-BP092424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022109.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.646	109	112	118	rBV	21452	32341	1.07%	0.196%
2	5.758	466	471	477	rVB	33227	48180	1.59%	0.292%
3	6.364	565	574	577	rVV8	7950	17667	0.58%	0.107%
4	6.711	629	633	639	rVV4	13682	25627	0.84%	0.155%
5	6.769	639	643	646	rVV	16058	24176	0.80%	0.147%
6	6.811	646	650	654	rVV2	18014	28263	0.93%	0.171%
7	6.881	654	662	668	rVV4	47305	98590	3.25%	0.598%
8	6.934	668	671	676	rVV	14927	24775	0.82%	0.150%
9	7.040	684	689	695	rVV	24915	43514	1.43%	0.264%
10	7.105	695	700	702	rVV3	13788	20423	0.67%	0.124%
11	7.140	702	706	709	rVV	11323	18871	0.62%	0.114%
12	7.240	713	723	734	rVV3	36344	85599	2.82%	0.519%
13	7.334	734	739	749	rVV2	80366	160263	5.28%	0.972%
14	7.611	777	786	793	rBV2	5733	18669	0.62%	0.113%
15	7.699	793	801	808	rBV	299608	500540	16.50%	3.036%
16	7.764	808	812	818	rVB	32778	52298	1.72%	0.317%
17	7.816	818	821	828	rVB2	11901	19442	0.64%	0.118%
18	8.240	887	893	897	rVB4	9732	19432	0.64%	0.118%
19	8.334	904	909	911	rBV2	20420	33114	1.09%	0.201%
20	8.399	916	920	926	rVV2	34935	60319	1.99%	0.366%
21	8.458	926	930	933	rVV2	14836	22151	0.73%	0.134%
22	8.493	933	936	944	rVB3	10563	19654	0.65%	0.119%
23	8.687	965	969	976	rVB2	13288	22895	0.75%	0.139%
24	8.787	976	986	990	rBV4	16390	35022	1.15%	0.212%
25	8.840	990	995	1001	rVV3	25542	49185	1.62%	0.298%
26	8.916	1002	1008	1013	rVB	74949	115875	3.82%	0.703%
27	9.034	1018	1028	1035	rBV10	10450	30800	1.02%	0.187%
28	9.111	1035	1041	1047	rBV3	15781	29717	0.98%	0.180%
29	9.558	1110	1117	1122	rBV3	26792	57140	1.88%	0.347%
30	9.616	1122	1127	1130	rVV6	12573	26197	0.86%	0.159%
31	9.752	1143	1150	1156	rVV4	10756	24774	0.82%	0.150%
32	9.899	1167	1175	1179	rVV9	9254	26291	0.87%	0.159%
33	10.093	1204	1208	1216	rVV	36225	66729	2.20%	0.405%
34	10.463	1258	1271	1278	rBV	393965	739609	24.38%	4.487%
35	10.593	1286	1293	1306	rBV3	49022	132128	4.36%	0.802%
36	10.840	1329	1335	1343	rBV	37234	61819	2.04%	0.375%

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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_P\Methods\SFAM-EPA-BP092424.MA.M
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37	10.963	1347	1356	1363	rBV7	9506	26156	0.86%	0.159%
38	11.487	1440	1445	1450	rVV2	15296	25950	0.86%	0.157%
39	11.557	1450	1457	1468	rVB	76013	122248	4.03%	0.742%
40	11.728	1477	1486	1494	rVB2	29317	54604	1.80%	0.331%
41	11.887	1503	1513	1516	rBV3	31433	56694	1.87%	0.344%
42	11.922	1516	1519	1524	rVV2	25227	38763	1.28%	0.235%
43	12.022	1531	1536	1538	rVV2	17188	28927	0.95%	0.175%
44	12.046	1538	1540	1548	rVV6	18467	35554	1.17%	0.216%
45	12.128	1548	1554	1563	rVB2	52770	95760	3.16%	0.581%
46	12.340	1586	1590	1600	rVB6	10585	21690	0.72%	0.132%
47	12.540	1617	1624	1634	rVV9	13479	39308	1.30%	0.238%
48	12.722	1647	1655	1660	rVB5	9505	24150	0.80%	0.147%
49	13.051	1707	1711	1717	rVB2	16019	21927	0.72%	0.133%
50	13.140	1720	1726	1733	rVV2	36388	60873	2.01%	0.369%
51	13.269	1738	1748	1756	rBV3	19118	44475	1.47%	0.270%
52	13.357	1756	1763	1771	rVB3	8488	22306	0.74%	0.135%
53	13.493	1777	1786	1795	rBV8	14634	46709	1.54%	0.283%
54	13.575	1795	1800	1803	rBV	33827	47261	1.56%	0.287%
55	13.622	1803	1808	1812	rVV4	13212	28089	0.93%	0.170%
56	13.669	1813	1816	1818	rVV3	18066	26142	0.86%	0.159%
57	13.710	1818	1823	1829	rVV	99822	143092	4.72%	0.868%
58	13.804	1831	1839	1843	rBV6	12823	30950	1.02%	0.188%
59	13.940	1858	1862	1866	rBV	52602	69109	2.28%	0.419%
60	13.998	1866	1872	1877	rVV	81580	134014	4.42%	0.813%
61	14.222	1905	1910	1915	rVB	189099	253385	8.35%	1.537%
62	14.310	1915	1925	1932	rBV	658935	1004620	33.12%	6.094%
63	14.387	1934	1938	1943	rVB7	15035	23446	0.77%	0.142%
64	14.463	1945	1951	1958	rVB	119816	189859	6.26%	1.152%
65	14.546	1958	1965	1971	rBV4	25306	61241	2.02%	0.372%
66	14.698	1988	1991	1999	rVB8	20438	42371	1.40%	0.257%
67	14.857	2014	2018	2021	rBV5	11879	19063	0.63%	0.116%
68	14.893	2021	2024	2027	rBV2	22241	29714	0.98%	0.180%
69	14.951	2029	2034	2040	rVV	106680	163947	5.40%	0.995%
70	15.004	2040	2043	2049	rVB2	44413	63877	2.11%	0.388%
71	15.081	2051	2056	2062	rBV	132855	209465	6.90%	1.271%
72	15.187	2067	2074	2079	rVB4	49759	88763	2.93%	0.538%
73	15.251	2079	2085	2089	rBV7	11740	22327	0.74%	0.135%
74	15.316	2089	2096	2103	rVB	158075	221405	7.30%	1.343%
75	15.434	2109	2116	2124	rBV2	41234	75084	2.48%	0.455%
76	15.598	2139	2144	2149	rBV3	19000	31780	1.05%	0.193%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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77	15.687	2155	2159	2165	rVB2	34464	53593	1.77%	0.325%
78	15.969	2203	2207	2213	rVV	56587	77296	2.55%	0.469%
79	16.069	2219	2224	2233	rVB2	47853	89362	2.95%	0.542%
80	16.281	2257	2260	2264	rVV2	22133	26063	0.86%	0.158%
81	16.416	2278	2283	2287	rBV2	25525	37386	1.23%	0.227%
82	16.751	2333	2340	2354	rBV	2067439	3033535	100.00%	18.403%
83	16.881	2358	2362	2367	rVV	45304	70370	2.32%	0.427%
84	16.940	2367	2372	2382	rVB4	26659	55104	1.82%	0.334%
85	17.104	2394	2400	2410	rVV	951793	1383317	45.60%	8.392%
86	17.216	2413	2419	2425	rVV	148251	280640	9.25%	1.702%
87	17.269	2425	2428	2432	rVB3	28796	35380	1.17%	0.215%
88	17.416	2448	2453	2462	rVB3	36391	66550	2.19%	0.404%
89	17.645	2487	2492	2497	rVB2	38173	52122	1.72%	0.316%
90	17.716	2500	2504	2506	rVV4	12828	19118	0.63%	0.116%
91	17.745	2506	2509	2517	rVV8	22442	34984	1.15%	0.212%
92	18.045	2557	2560	2568	rVB7	24477	39103	1.29%	0.237%
93	18.363	2610	2614	2619	rBV3	25979	36789	1.21%	0.223%
94	19.016	2721	2725	2728	rBV3	23230	37729	1.24%	0.229%
95	19.122	2740	2743	2748	rBV6	16476	19992	0.66%	0.121%
96	19.586	2817	2822	2827	rBV2	233610	331214	10.92%	2.009%
97	20.181	2920	2923	2928	rBV5	29975	47765	1.57%	0.290%
98	21.539	3148	3154	3162	rBV2	1217644	1854481	61.13%	11.250%
99	24.598	3670	3674	3683	rVB2	131607	289995	9.56%	1.759%
100	24.827	3705	3713	3726	rVB2	780101	1975181	65.11%	11.982%

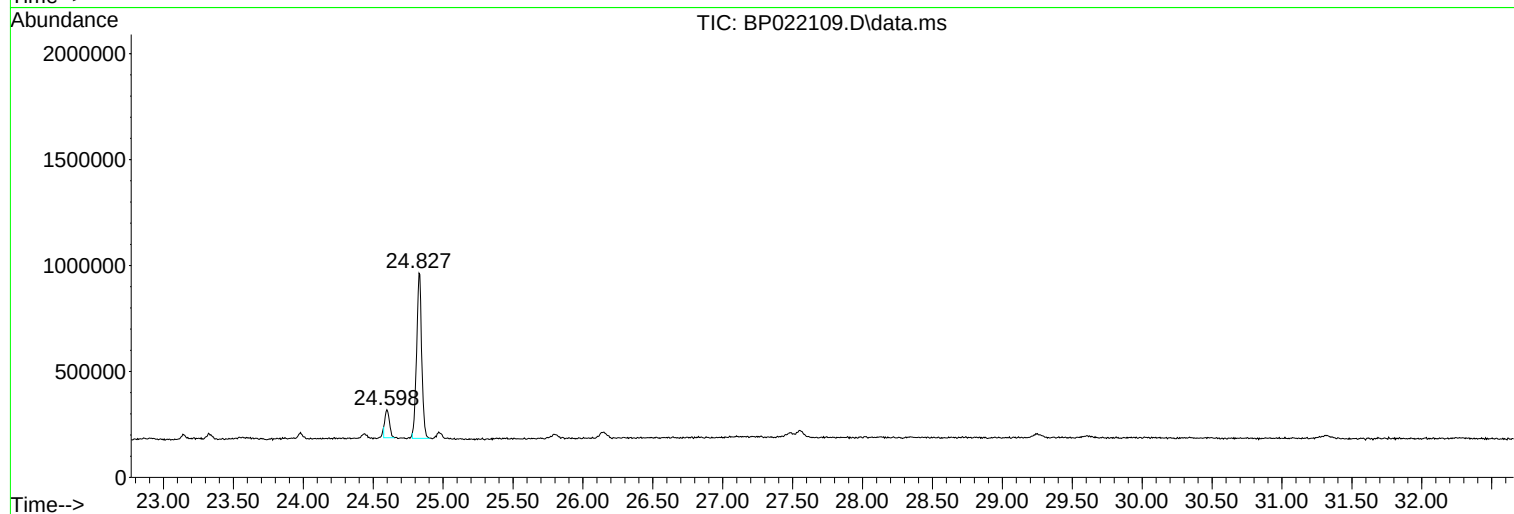
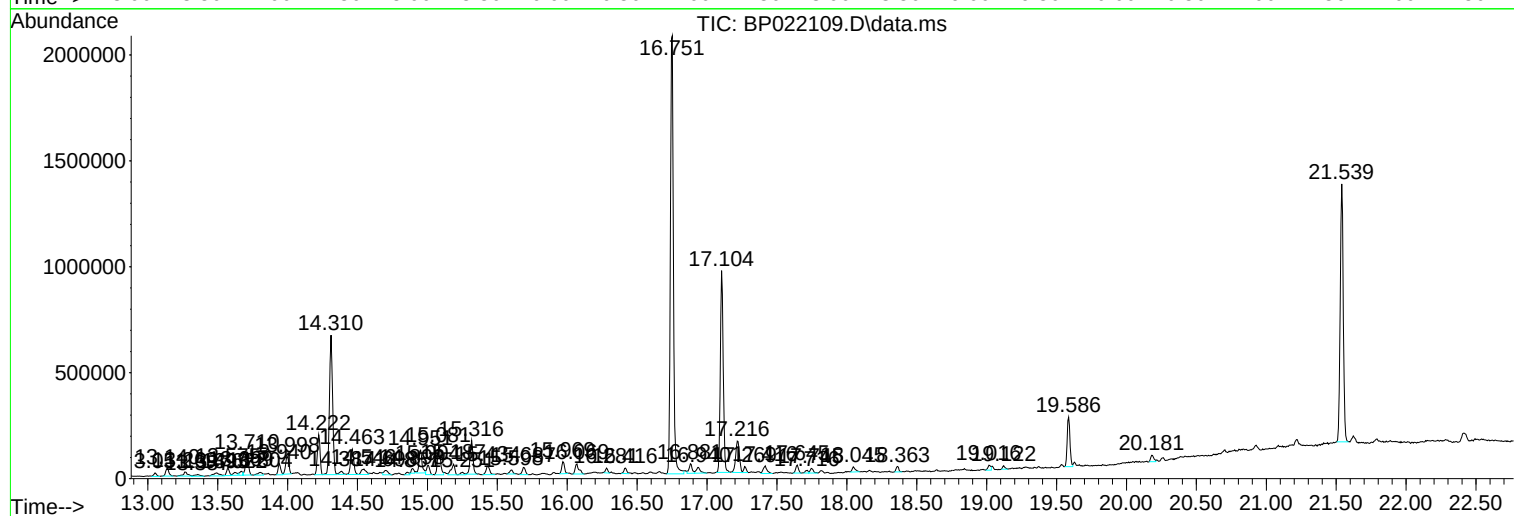
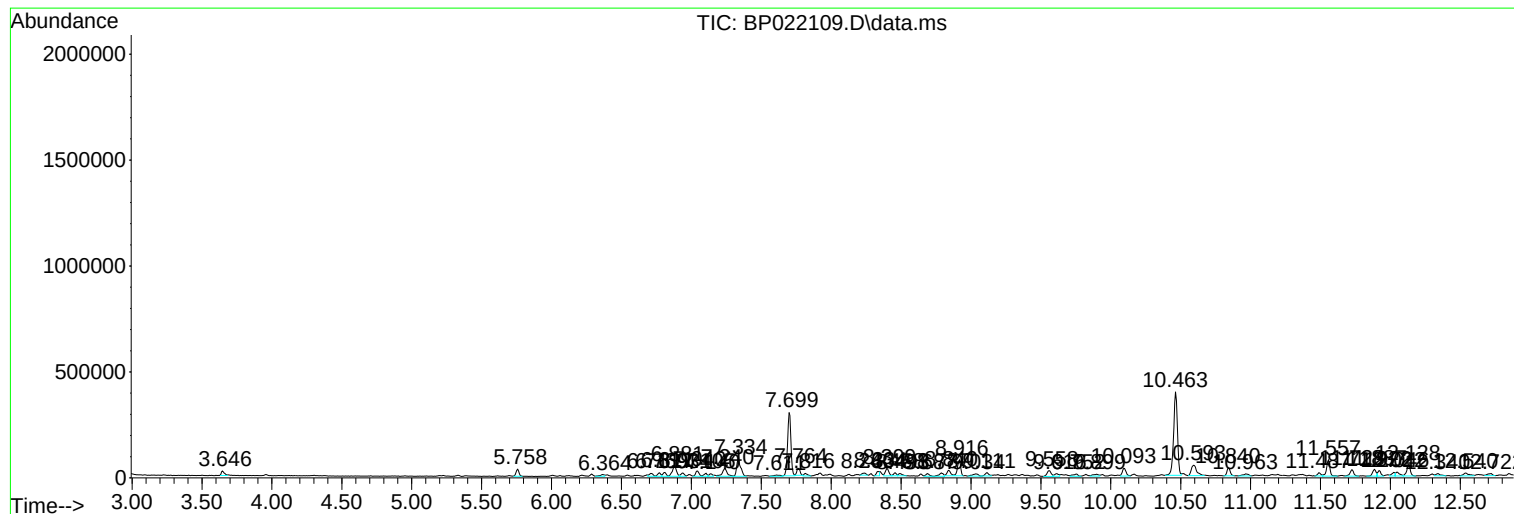
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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



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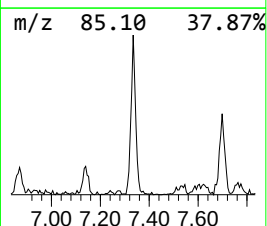
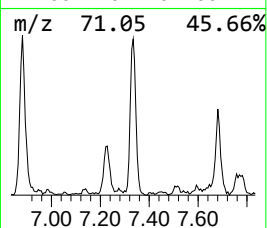
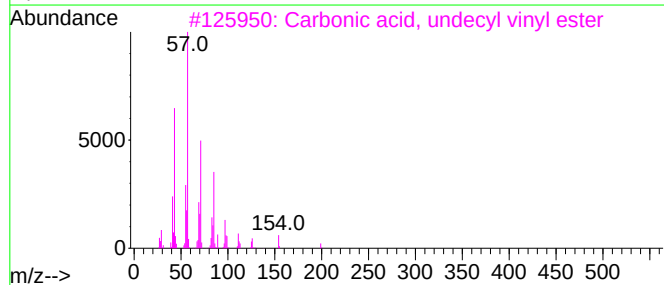
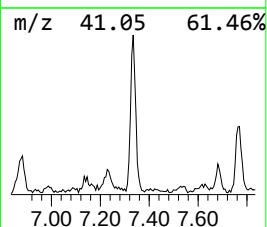
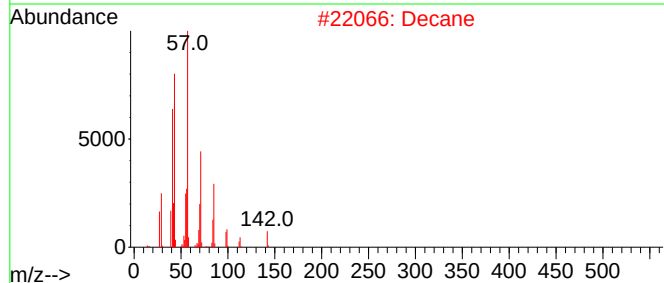
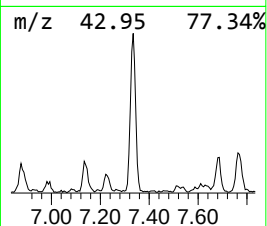
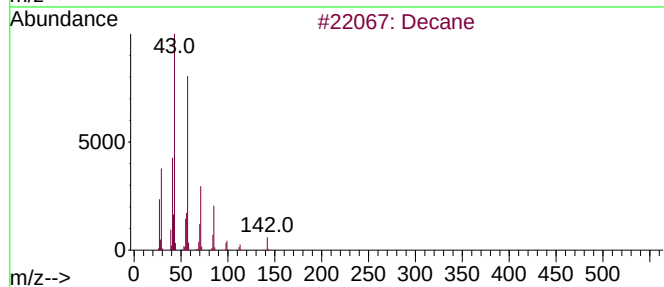
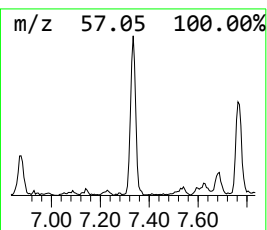
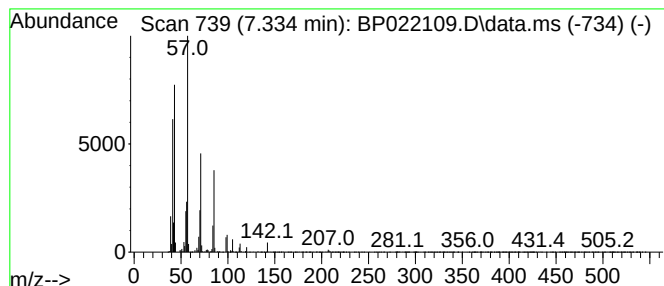
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.334	6.40 ng/ul	160263	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	95
2			Decane	142	C10H22	000124-18-5	91
3			Carbonic acid, undecyl vinyl ester	242	C14H26O3	1000382-54-9	83
4			Decane	142	C10H22	000124-18-5	83
5			Decane, 1,1'-oxybis-	298	C20H42O	002456-28-2	80



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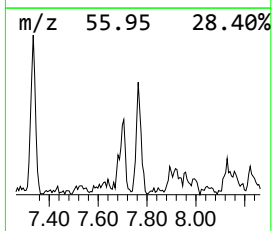
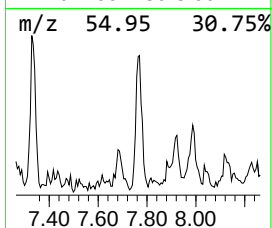
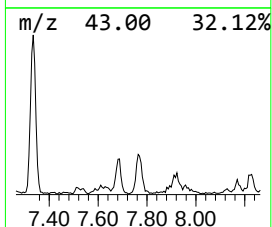
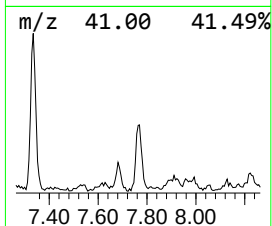
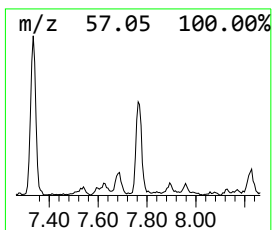
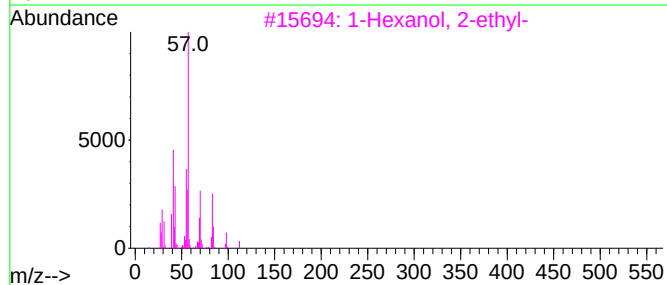
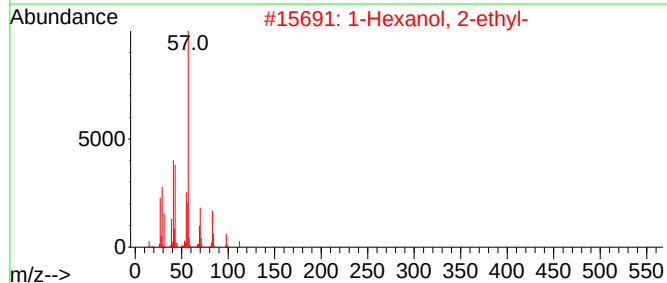
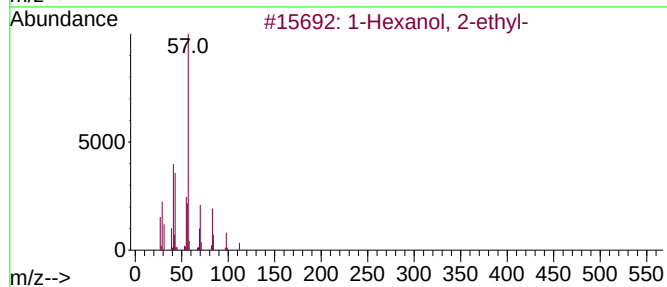
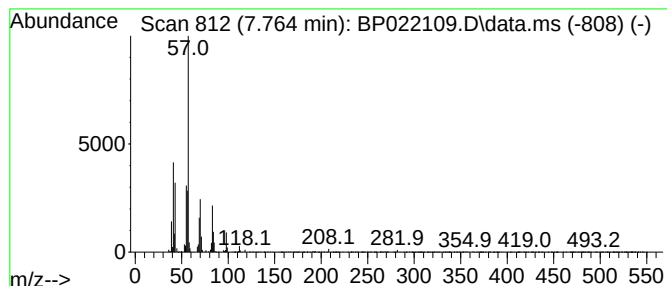
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.764	2.09 ng/ul	52298	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
4	Cyclopropane, 1-(2-methylbutyl)-...			168	C12H24	064723-36-0	58
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	53



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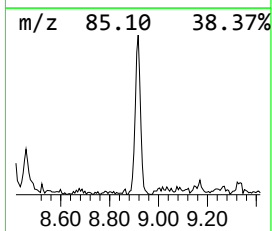
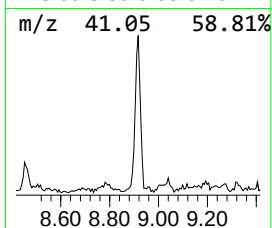
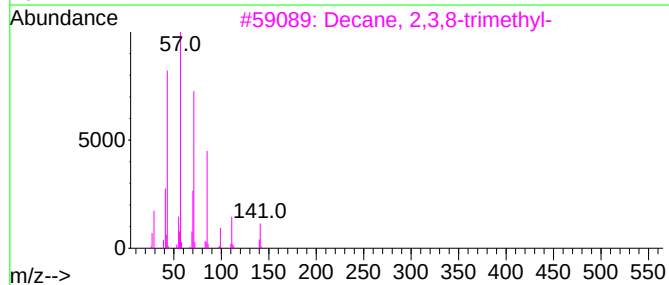
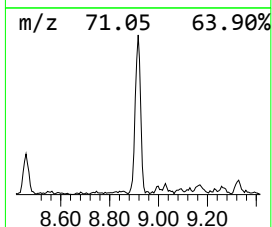
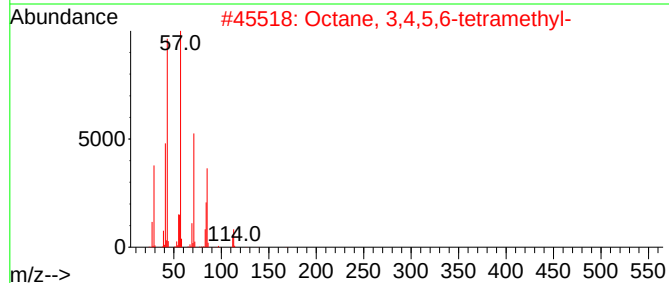
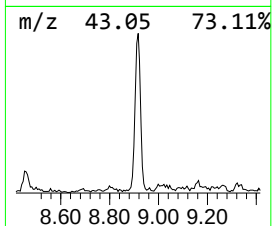
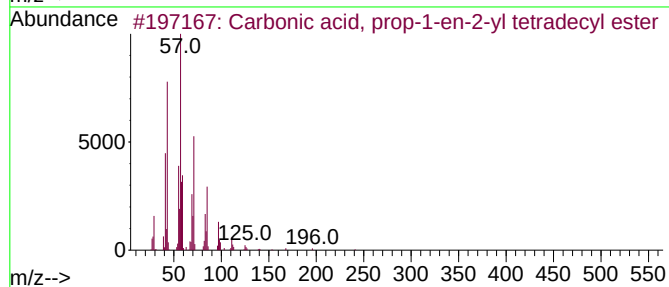
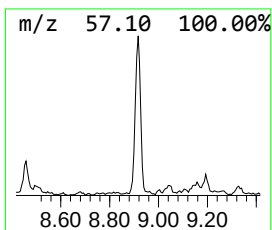
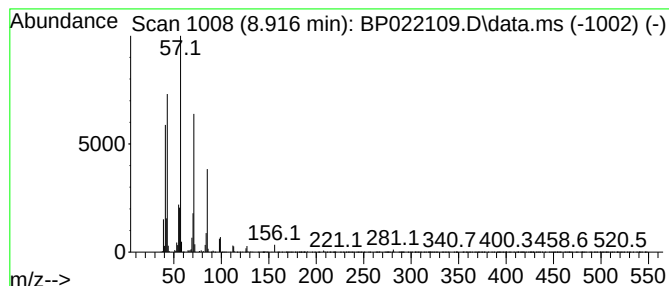
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Carbonic acid, prop-1-en-2-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.916	4.63 ng/ul	115875	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	86
2			Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	72
3			Decane, 2,3,8-trimethyl-	184	C13H28	062238-14-6	64
4			Undecane	156	C11H24	001120-21-4	58
5			Tridecane	184	C13H28	000629-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022109.D
Acq On : 28 Sep 2024 17:48
Operator : RC/JU
Sample : P3910-10MEDL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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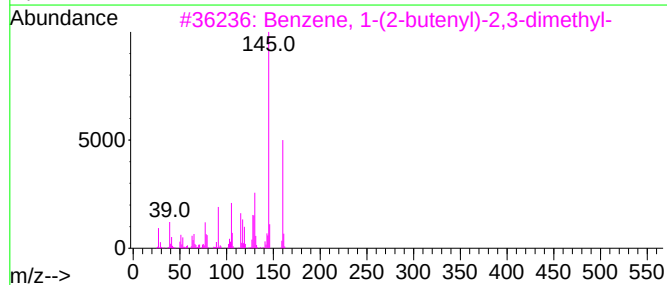
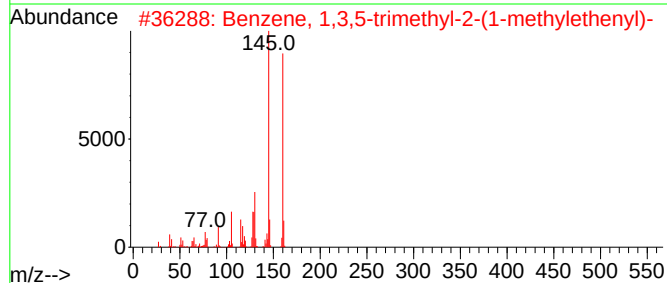
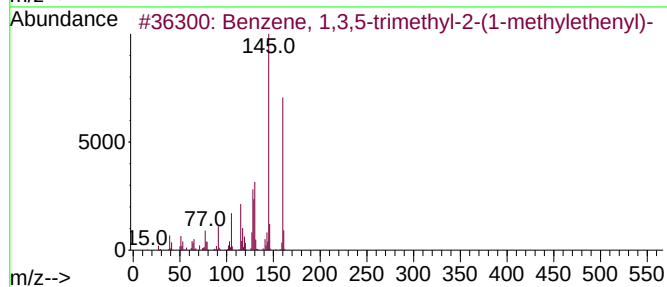
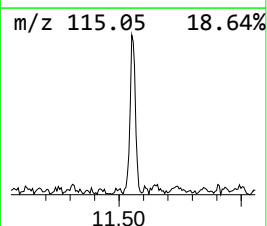
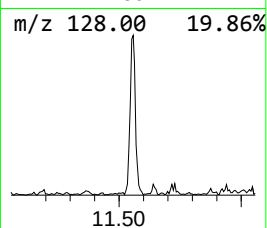
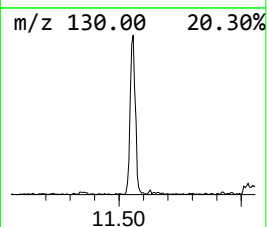
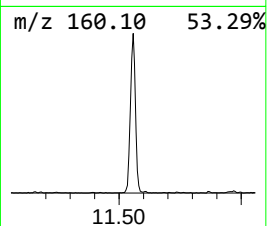
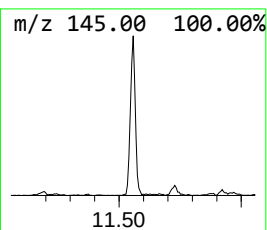
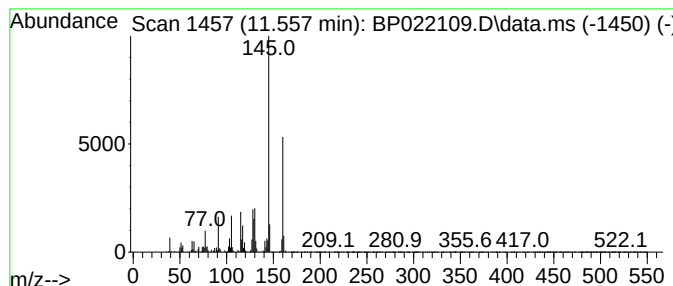
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.557	3.31 ng/ul	122248	Naphthalene-d8	10.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	96
2			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	94
3			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
4			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	94
5			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022109.D
Acq On : 28 Sep 2024 17:48
Operator : RC/JU
Sample : P3910-10MEDL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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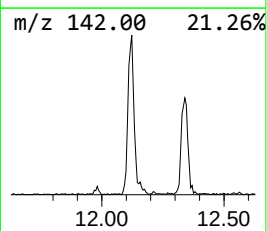
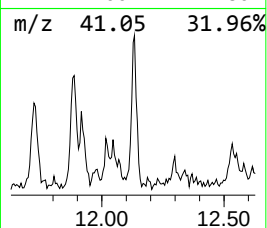
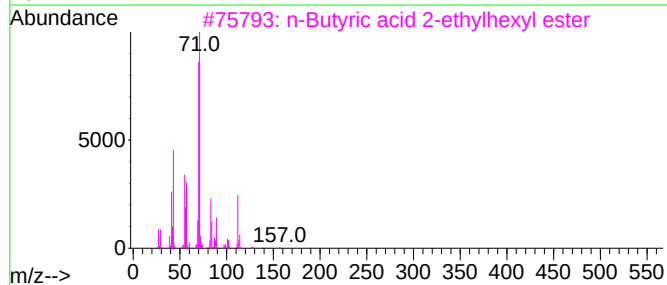
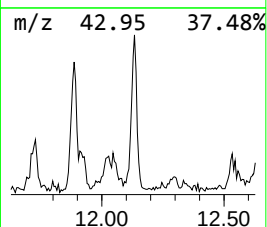
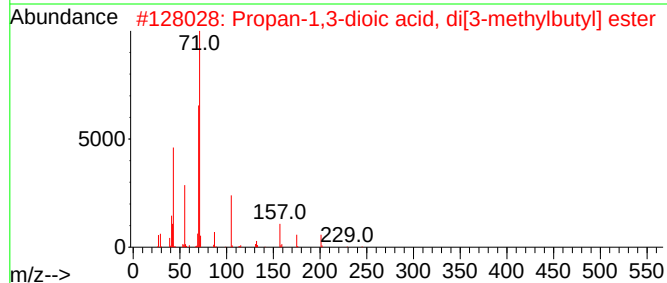
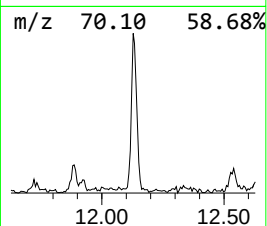
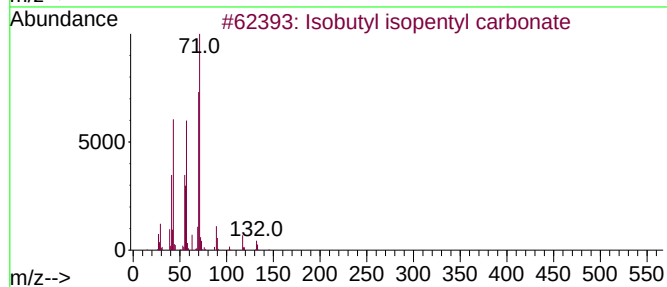
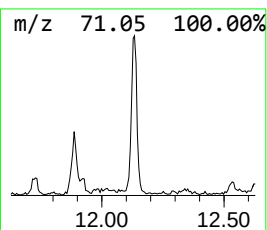
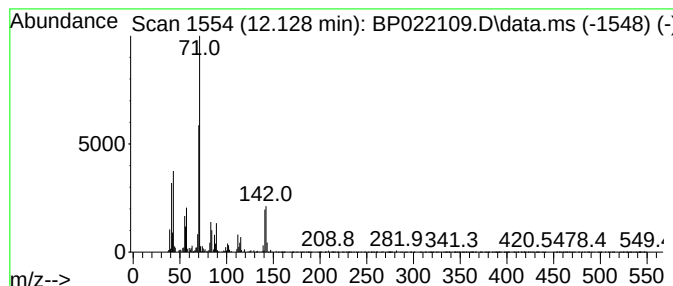
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Isobutyl isopentyl carbonate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.128	2.59 ng/ul	95760	Naphthalene-d8	10.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl isopentyl carbonate	188	C10H20O3	1000372-76-3	50
2			Propan-1,3-dioic acid, di[3-meth...	244	C13H24O4	064617-96-5	47
3			n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	46
4			2-Nonanol, 2-methylpropionate	214	C13H26O2	069121-76-2	43
5			Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022109.D
Acq On : 28 Sep 2024 17:48
Operator : RC/JU
Sample : P3910-10MEDL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

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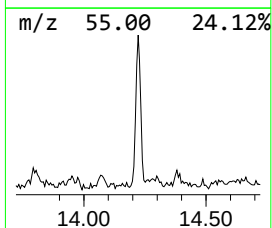
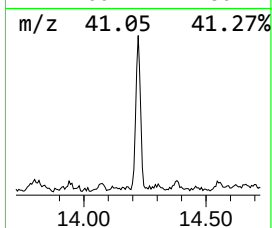
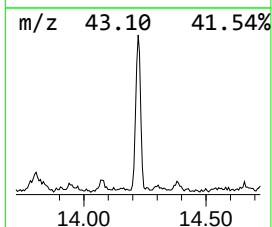
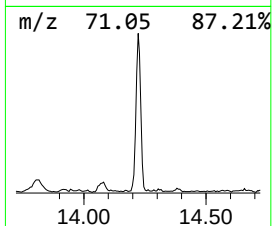
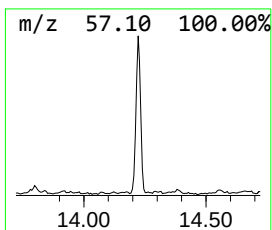
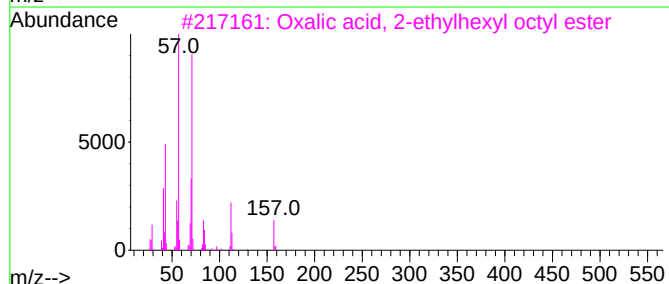
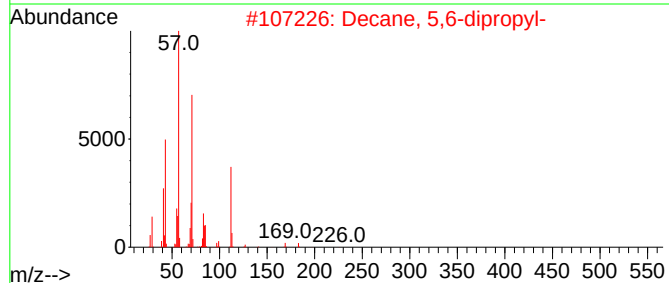
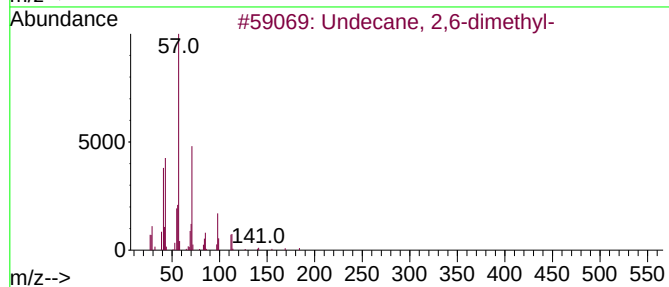
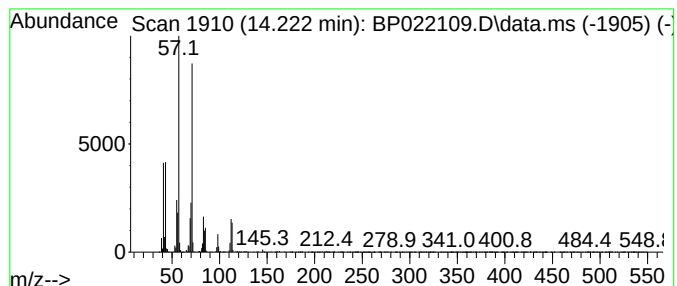
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.222	5.04 ng/ul	253385	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	80
2			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	80
3			Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	72
4			Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
5			2-Undecene, 5-methyl-	168	C12H24	056851-34-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022109.D
Acq On : 28 Sep 2024 17:48
Operator : RC/JU
Sample : P3910-10MEDL 10X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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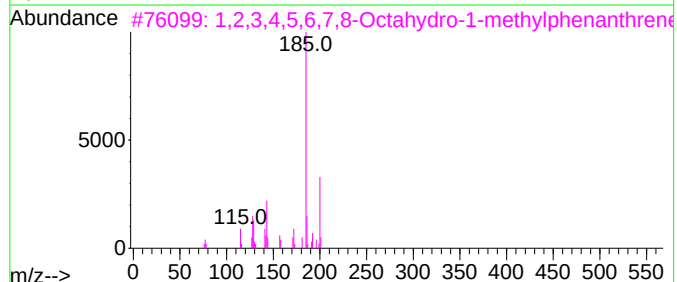
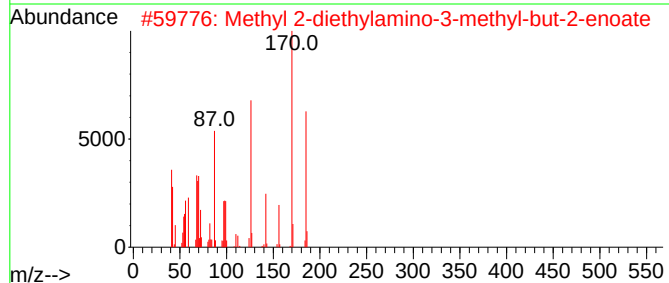
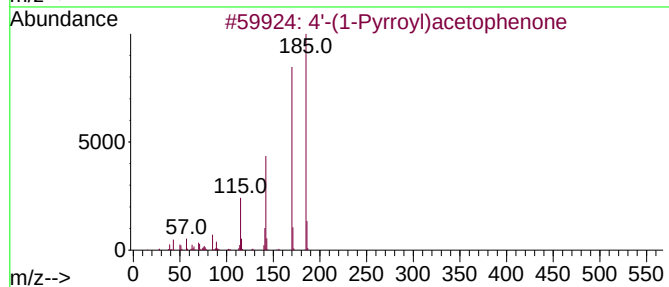
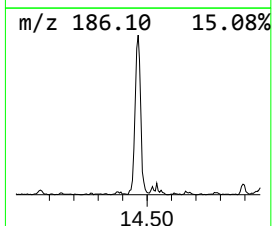
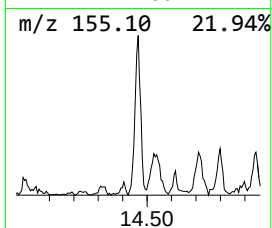
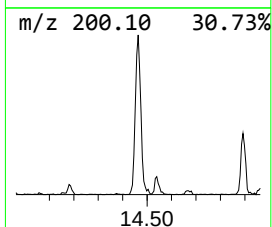
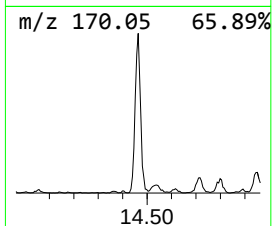
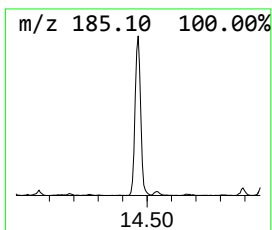
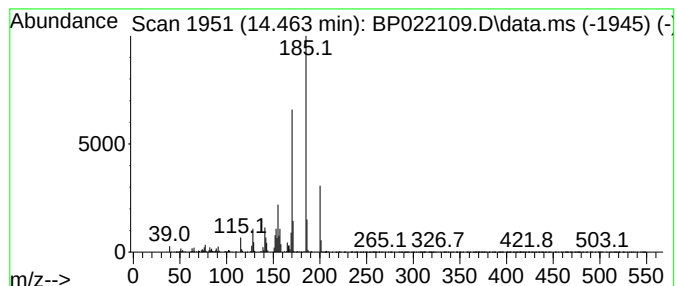
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 4'-(1-Pyrrolyl)acetophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	3.78 ng/ul	189859	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	52		
2	Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	47		
3	1,2,3,4,5,6,7,8-Octahydro-1-meth...	200	C15H20	1000080-19-4	46		
4	1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46		
5	4-Chlorophenol, TMS derivative	200	C9H13ClOSi	017005-59-3	46		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022109.D
Acq On : 28 Sep 2024 17:48
Operator : RC/JU
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Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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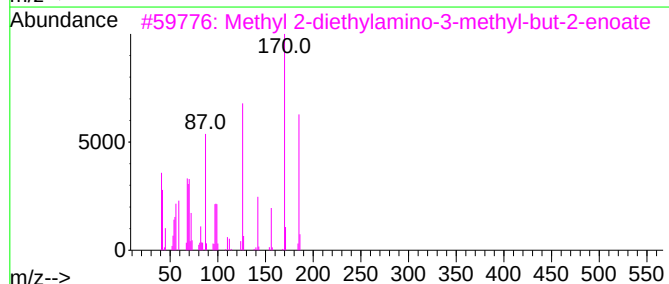
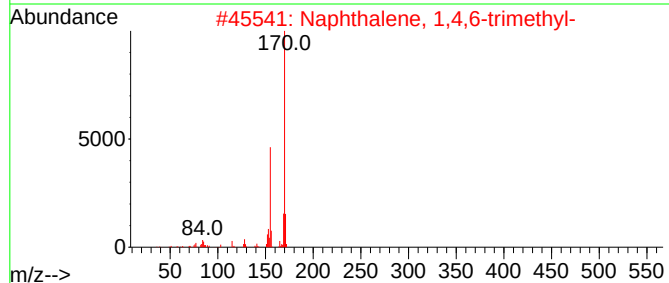
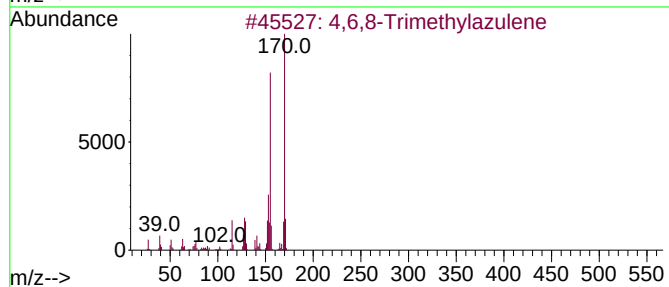
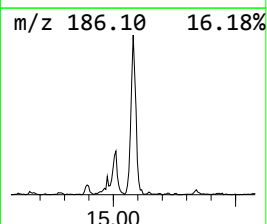
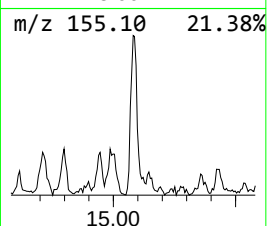
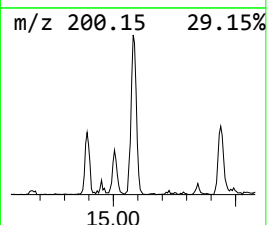
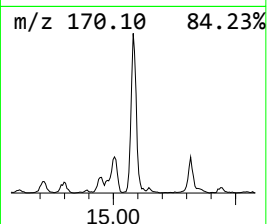
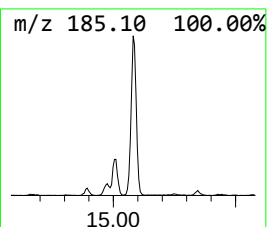
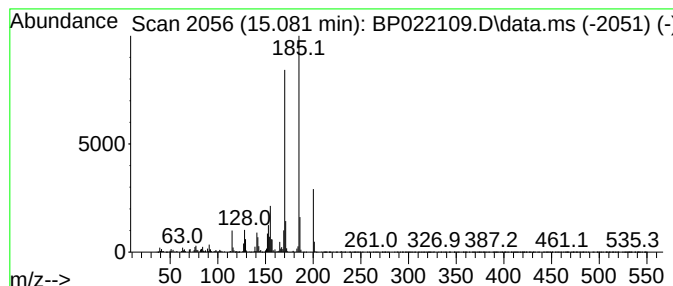
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.081	4.17 ng/ul	209465	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	50
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
3			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	50
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	46
5			Methyl 2-(4-methyl-6-phenylcyclo...	244	C16H20O2	1000215-33-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022109.D
Acq On : 28 Sep 2024 17:48
Operator : RC/JU
Sample : P3910-10MEDL 10X
Misc :
ALS Vial : 15 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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1-Hexanol, 2-et...	7.764	2.1	ng/ul	52298	1	7.699	500540	20.0
Carbonic acid, ...	8.916	4.6	ng/ul	115875	1	7.699	500540	20.0
Benzene, 1,3,5-...	11.557	3.3	ng/ul	122248	2	10.463	739609	20.0
Isobutyl isopen...	12.128	2.6	ng/ul	95760	2	10.463	739609	20.0
(DEL) Alkane: S...	14.222	5.0	ng/ul	253385	3	14.310	1004620	20.0
4'-(1-Pyrroyl)a...	14.463	3.8	ng/ul	189859	3	14.310	1004620	20.0
4,6,8-Trimethyl...	15.081	4.2	ng/ul	209465	3	14.310	1004620	20.0