

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
 Data File : BP022108.D
 Acq On : 28 Sep 2024 17:07
 Operator : RC/JU
 Sample : P3910-08MEDL 20X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DDC18MEDL

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: BP022108.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.652	109	113	123	rBV	11037	19465	0.78%	0.137%
2	5.758	464	471	476	rBV	29439	42534	1.70%	0.299%
3	6.711	630	633	638	rVV4	12149	19910	0.80%	0.140%
4	6.770	638	643	646	rVV2	14706	21033	0.84%	0.148%
5	6.811	646	650	654	rVV2	15064	24181	0.97%	0.170%
6	6.875	654	661	668	rVV4	29290	67459	2.70%	0.474%
7	6.934	668	671	676	rVV	14279	21589	0.86%	0.152%
8	7.040	684	689	694	rVV2	13996	23853	0.95%	0.167%
9	7.099	694	699	702	rVV2	12454	19425	0.78%	0.136%
10	7.140	702	706	712	rVV4	10544	21211	0.85%	0.149%
11	7.240	712	723	730	rVV4	21675	56767	2.27%	0.399%
12	7.334	733	739	752	rVV2	75411	144383	5.77%	1.014%
13	7.699	793	801	808	rVV	324877	535773	21.43%	3.762%
14	7.764	808	812	817	rVV3	31516	56575	2.26%	0.397%
15	7.817	817	821	827	rVB3	10561	19344	0.77%	0.136%
16	7.917	835	838	843	rVV2	13013	21495	0.86%	0.151%
17	8.187	878	884	887	rVV6	7134	15517	0.62%	0.109%
18	8.228	887	891	897	rVV3	13949	30972	1.24%	0.217%
19	8.334	904	909	916	rVV5	16913	45680	1.83%	0.321%
20	8.405	916	921	926	rVV2	15802	29681	1.19%	0.208%
21	8.458	926	930	933	rVV2	11963	19490	0.78%	0.137%
22	8.493	933	936	944	rVB2	9445	16714	0.67%	0.117%
23	8.640	955	961	965	rBV	9482	15359	0.61%	0.108%
24	8.687	965	969	974	rVB	9757	15775	0.63%	0.111%
25	8.787	978	986	990	rBV5	12778	26732	1.07%	0.188%
26	8.840	990	995	999	rVV2	16596	27288	1.09%	0.192%
27	8.911	1002	1007	1013	rVB	64638	101342	4.05%	0.712%
28	9.034	1018	1028	1035	rBV	9064	25087	1.00%	0.176%
29	9.111	1035	1041	1047	rBV4	15337	27724	1.11%	0.195%
30	9.363	1081	1084	1093	rVB4	7058	16162	0.65%	0.113%
31	9.552	1107	1116	1121	rBV7	16775	36245	1.45%	0.254%
32	9.893	1165	1174	1178	rBV9	6568	16243	0.65%	0.114%
33	10.093	1203	1208	1214	rVB3	15824	29558	1.18%	0.208%
34	10.458	1261	1270	1277	rVV	425015	755558	30.22%	5.305%
35	10.581	1285	1291	1306	rBV6	31016	92740	3.71%	0.651%
36	10.834	1329	1334	1342	rVB2	31734	53728	2.15%	0.377%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
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37	10.969	1353	1357	1363	rVB5	8944	15816	0.63%	0.111%
38	11.358	1415	1423	1424	rVV8	6055	14575	0.58%	0.102%
39	11.481	1439	1444	1449	rVV4	13031	24926	1.00%	0.175%
40	11.552	1449	1456	1463	rVB	58024	101500	4.06%	0.713%
41	11.716	1478	1484	1493	rVB4	26851	45242	1.81%	0.318%
42	11.881	1503	1512	1515	rBV2	29774	49224	1.97%	0.346%
43	11.922	1515	1519	1524	rVV2	22165	35640	1.43%	0.250%
44	12.016	1532	1535	1537	rVV2	15037	22547	0.90%	0.158%
45	12.046	1537	1540	1547	rVV4	15118	32144	1.29%	0.226%
46	12.128	1547	1554	1562	rVB3	47264	85452	3.42%	0.600%
47	12.528	1614	1622	1627	rBV10	11883	24741	0.99%	0.174%
48	12.710	1646	1653	1659	rVB5	10380	24934	1.00%	0.175%
49	12.846	1671	1676	1685	rBV8	8575	17722	0.71%	0.124%
50	13.052	1703	1711	1716	rVV2	12668	23668	0.95%	0.166%
51	13.134	1719	1725	1730	rBV	30403	50207	2.01%	0.353%
52	13.263	1743	1747	1755	rVV2	16946	34729	1.39%	0.244%
53	13.352	1757	1762	1768	rVB8	6223	14244	0.57%	0.100%
54	13.569	1794	1799	1805	rVV2	26789	42379	1.69%	0.298%
55	13.628	1805	1809	1811	rVV	12364	20236	0.81%	0.142%
56	13.663	1811	1815	1816	rVV4	13110	21520	0.86%	0.151%
57	13.710	1819	1823	1828	rVB	46509	69049	2.76%	0.485%
58	13.799	1834	1838	1843	rVB5	9502	16499	0.66%	0.116%
59	13.946	1857	1863	1867	rBV	42165	61528	2.46%	0.432%
60	13.999	1867	1872	1877	rVV	46219	72963	2.92%	0.512%
61	14.057	1879	1882	1889	rVB4	13339	26002	1.04%	0.183%
62	14.222	1905	1910	1915	rBV	172907	213111	8.52%	1.496%
63	14.310	1918	1925	1931	rVV	794786	1040855	41.63%	7.308%
64	14.381	1933	1937	1942	rVV6	16016	26668	1.07%	0.187%
65	14.451	1944	1949	1957	rVB	108115	158255	6.33%	1.111%
66	14.528	1957	1962	1970	rBV5	14688	36206	1.45%	0.254%
67	14.704	1988	1992	1998	rVB6	23248	40044	1.60%	0.281%
68	14.851	2012	2017	2020	rBV7	8819	16454	0.66%	0.116%
69	14.893	2020	2024	2026	rBV3	17260	22799	0.91%	0.160%
70	14.951	2029	2034	2041	rVV	80546	126494	5.06%	0.888%
71	15.010	2041	2044	2050	rVB3	34017	48387	1.94%	0.340%
72	15.081	2050	2056	2061	rBV	112775	171387	6.85%	1.203%
73	15.187	2072	2074	2079	rVB	33695	44401	1.78%	0.312%
74	15.322	2092	2097	2103	rVB	73628	107078	4.28%	0.752%
75	15.428	2110	2115	2123	rBV2	26357	44144	1.77%	0.310%
76	15.604	2140	2145	2149	rVV2	15356	24176	0.97%	0.170%

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77	15.693	2155	2160	2166	rVB6	26105	40433	1.62%	0.284%
78	15.904	2192	2196	2199	rBV6	9422	14427	0.58%	0.101%
79	15.975	2203	2208	2214	rBV2	38904	58228	2.33%	0.409%
80	16.063	2218	2223	2226	rBV2	39760	61841	2.47%	0.434%
81	16.287	2257	2261	2266	rBV	16459	19364	0.77%	0.136%
82	16.410	2278	2282	2288	rVB2	21564	29984	1.20%	0.211%
83	16.751	2334	2340	2355	rBV	1977471	2500500	100.00%	17.556%
84	16.881	2358	2362	2367	rVV3	33704	47361	1.89%	0.333%
85	16.928	2367	2370	2375	rVB6	18374	28443	1.14%	0.200%
86	17.104	2395	2400	2414	rVV	902501	1395554	55.81%	9.798%
87	17.216	2414	2419	2425	rVV	88563	158331	6.33%	1.112%
88	17.269	2425	2428	2433	rVB3	19714	30557	1.22%	0.215%
89	17.416	2449	2453	2461	rVB4	27516	49074	1.96%	0.345%
90	17.645	2487	2492	2498	rBV2	30911	46844	1.87%	0.329%
91	17.716	2500	2504	2508	rVV6	10748	16601	0.66%	0.117%
92	17.757	2508	2511	2517	rVB8	14535	21065	0.84%	0.148%
93	18.051	2557	2561	2569	rVB7	17458	34489	1.38%	0.242%
94	18.363	2611	2614	2618	rBV	21089	26512	1.06%	0.186%
95	19.034	2724	2728	2730	rBV5	17327	26103	1.04%	0.183%
96	19.128	2741	2744	2749	rBV4	16184	18935	0.76%	0.133%
97	19.592	2818	2823	2827	rBV	122515	155181	6.21%	1.090%
98	20.192	2922	2925	2929	rBV4	21531	32511	1.30%	0.228%
99	21.545	3149	3155	3164	rVB	1163933	1841866	73.66%	12.932%
100	24.833	3706	3714	3726	rVB	771316	2008168	80.31%	14.099%

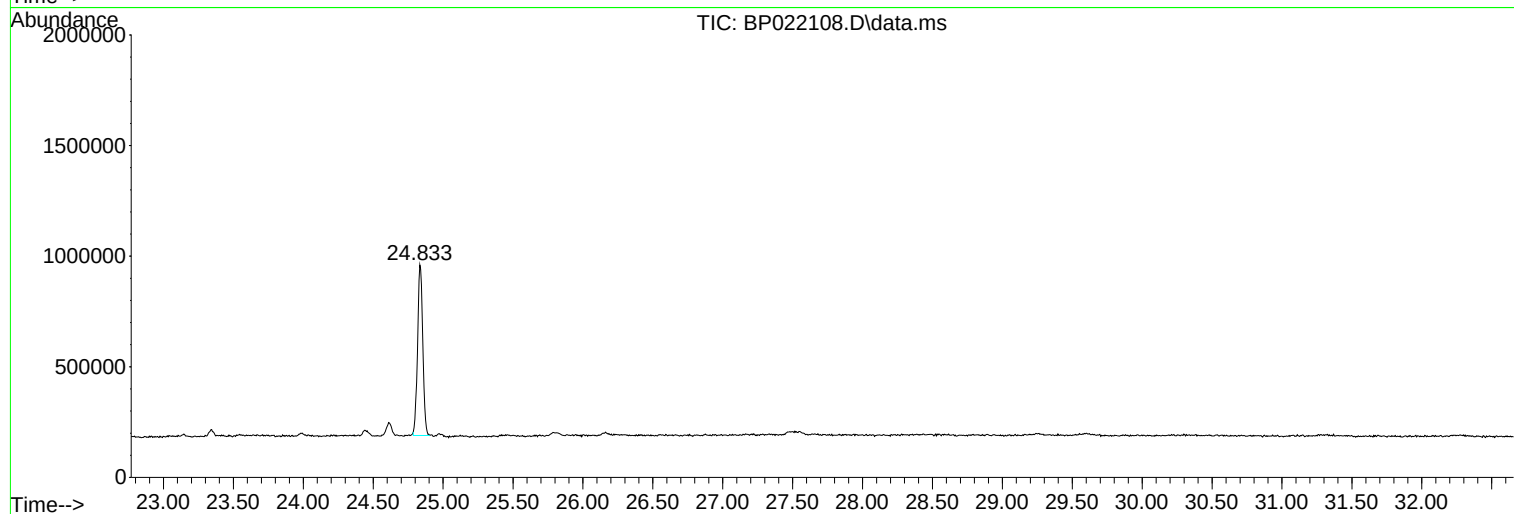
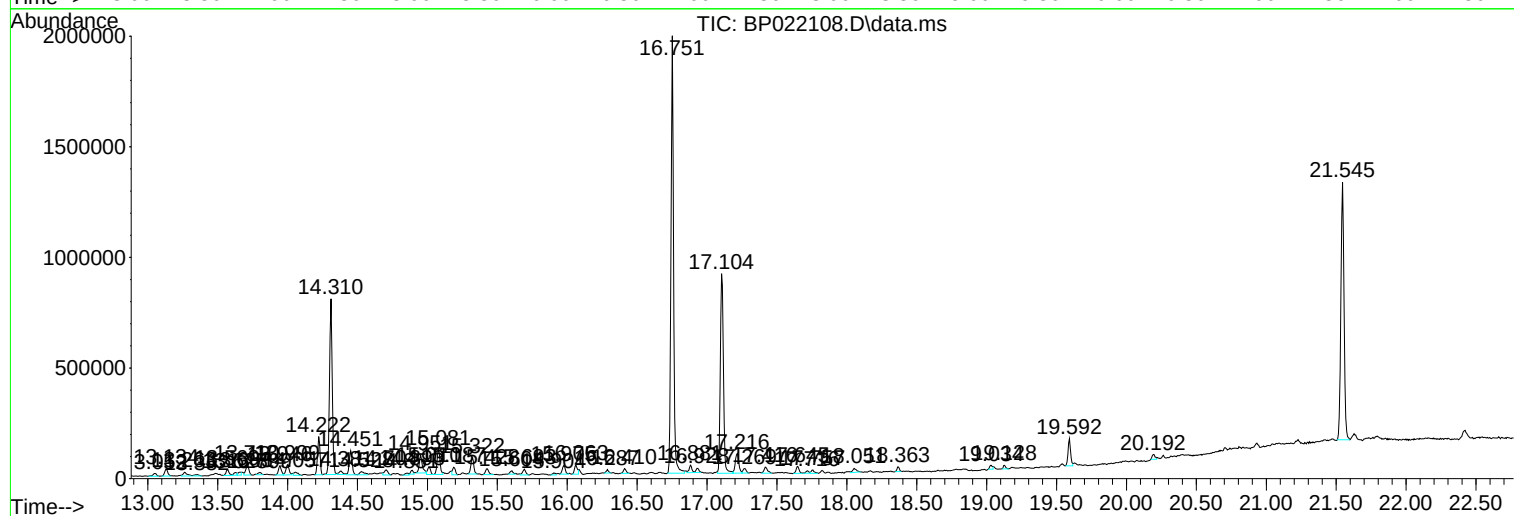
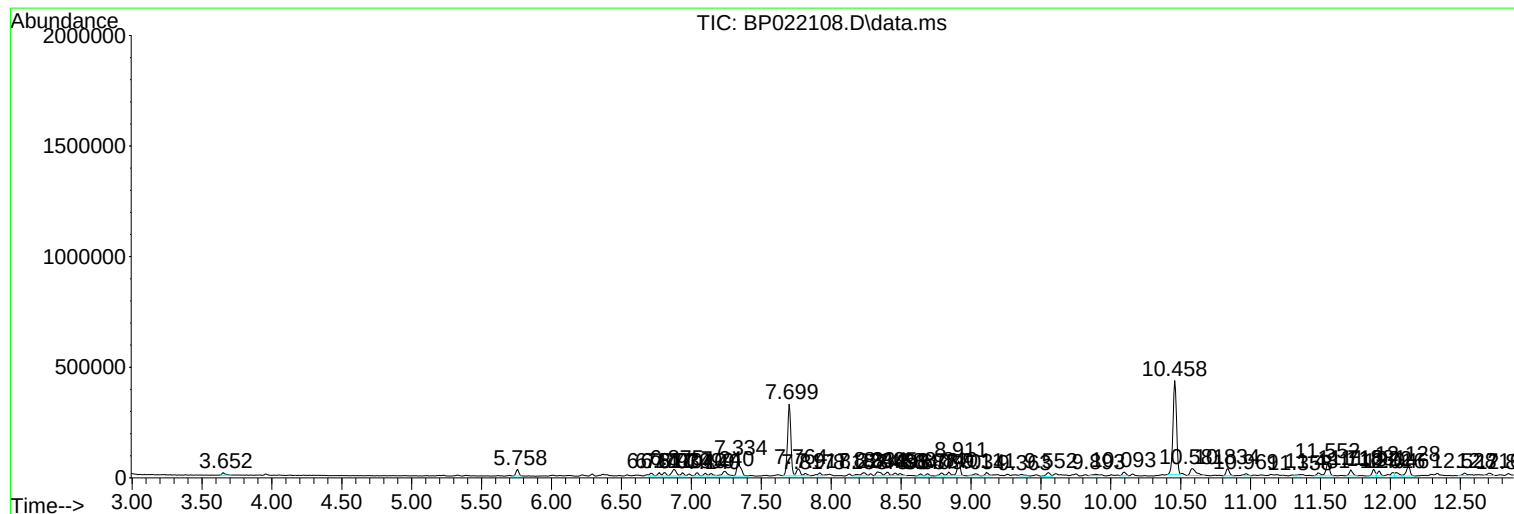
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ALS Vial : 14 Sample Multiplier: 1

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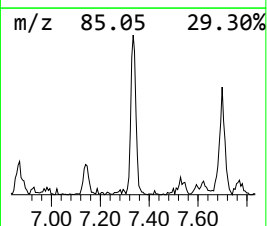
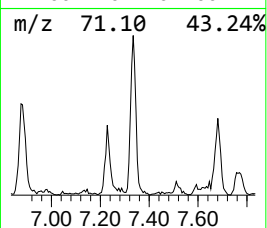
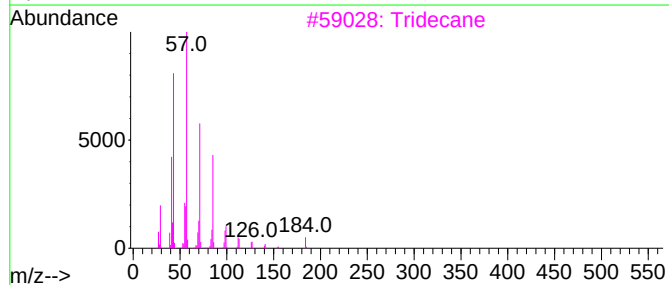
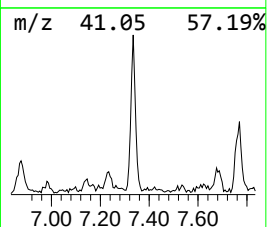
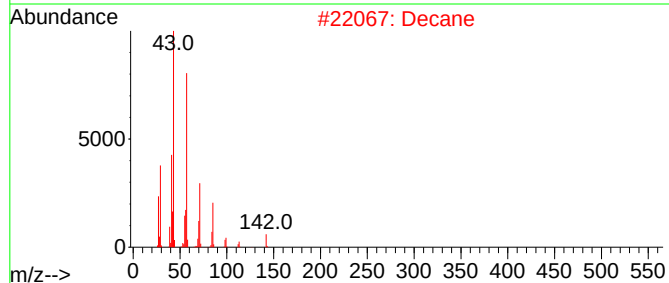
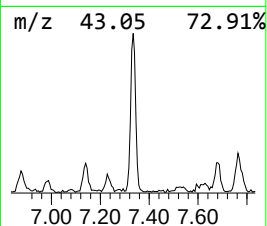
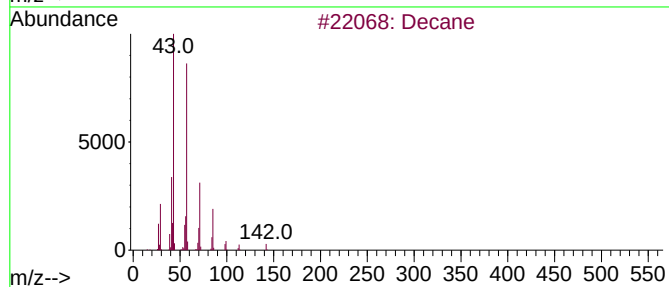
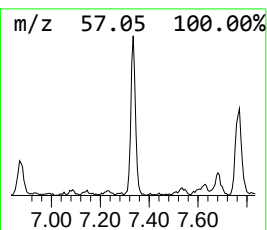
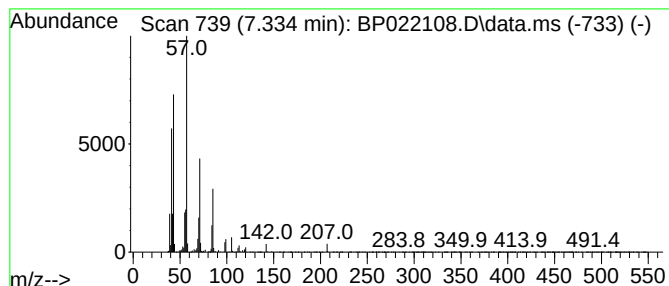
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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	5.39 ng/ul	144383	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	91
2			Decane	142	C10H22	000124-18-5	78
3			Tridecane	184	C13H28	000629-50-5	72
4			Hexadecane	226	C16H34	000544-76-3	64
5			Undecane	156	C11H24	001120-21-4	64



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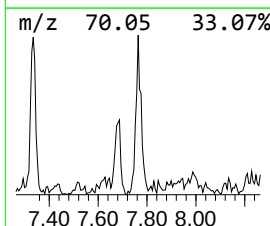
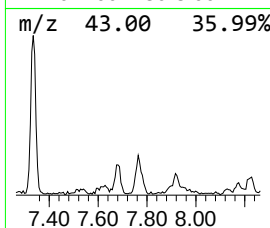
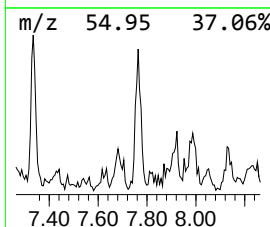
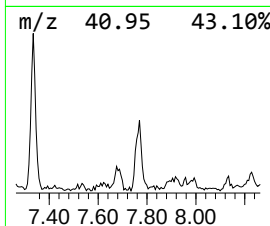
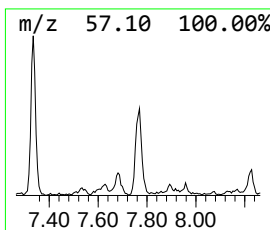
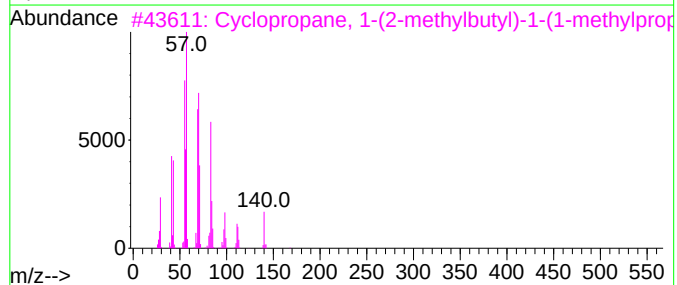
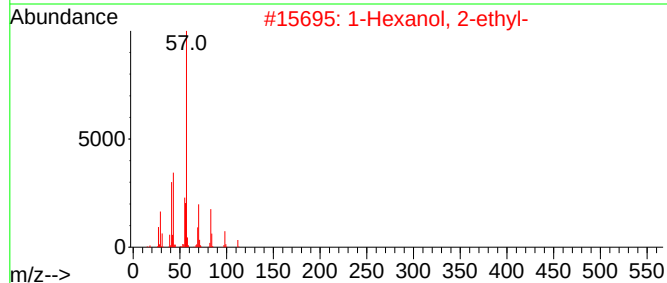
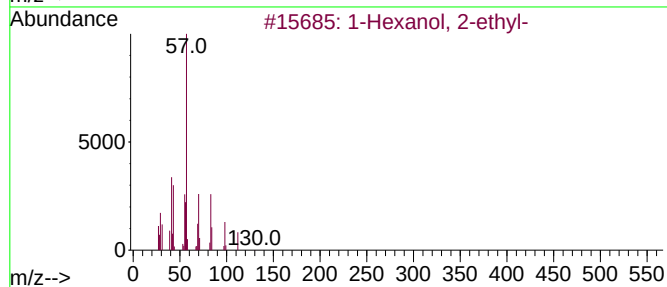
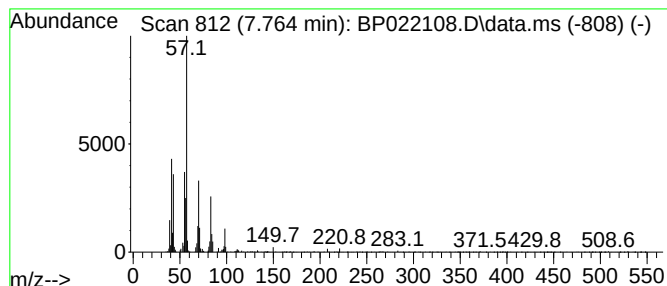
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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	2.11 ng/ul	56575	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	64
3	Cyclopropane, 1-(2-methylbutyl)-...			168	C12H24	064723-36-0	64
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
5	Carbonic acid, decyl 2-ethylhexy...			314	C19H38O3	1000383-13-7	50



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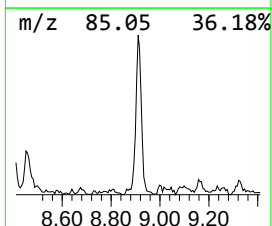
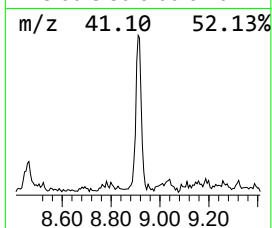
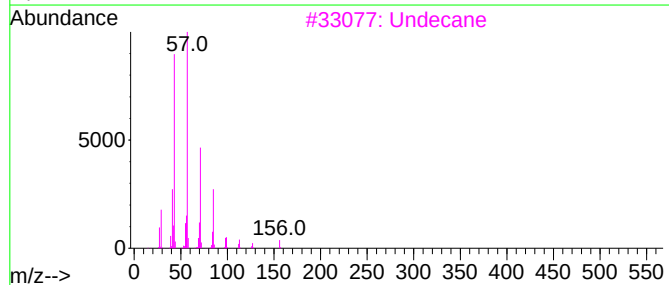
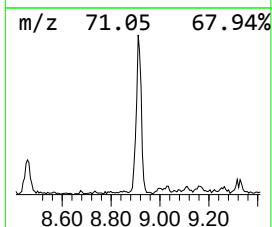
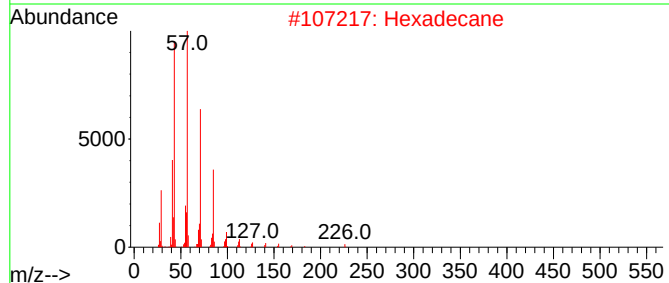
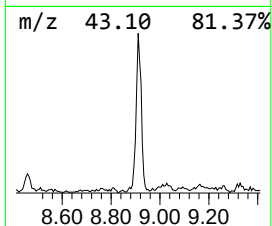
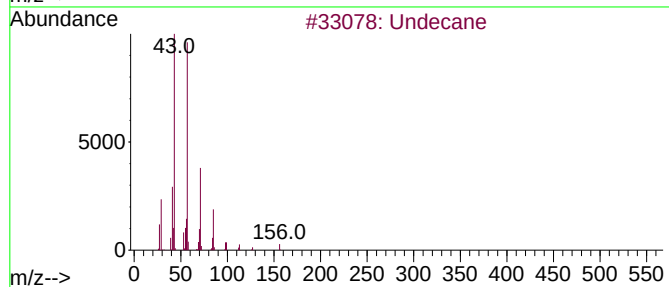
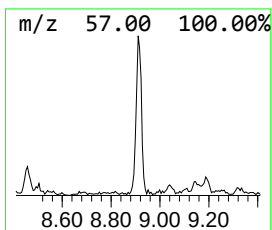
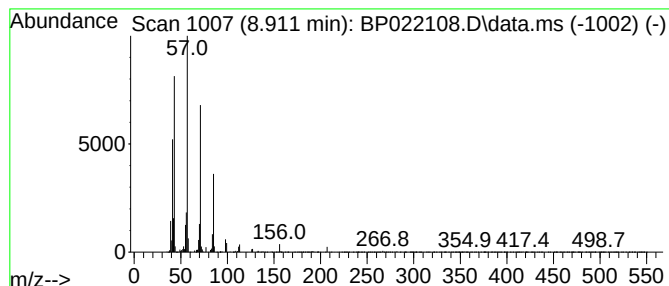
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.911	3.78 ng/ul	101342	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Hexadecane	226	C16H34	000544-76-3	90
3			Undecane	156	C11H24	001120-21-4	87
4			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	72
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022108.D
Acq On : 28 Sep 2024 17:07
Operator : RC/JU
Sample : P3910-08MEDL 20X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DDC18MEDL

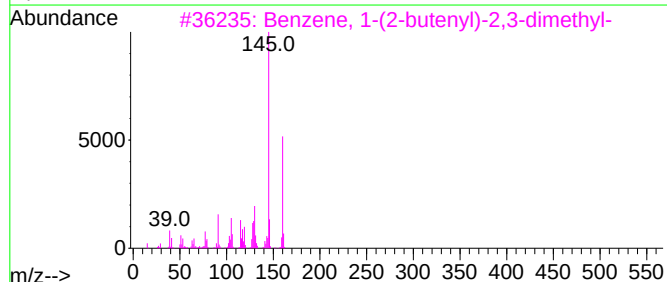
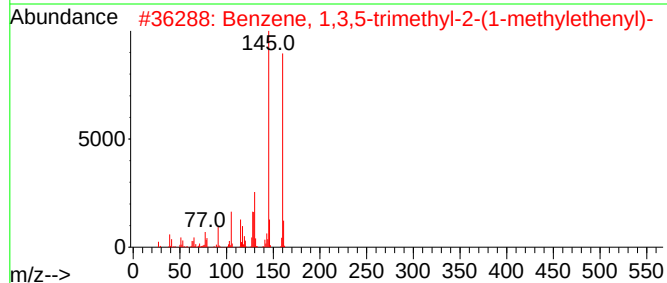
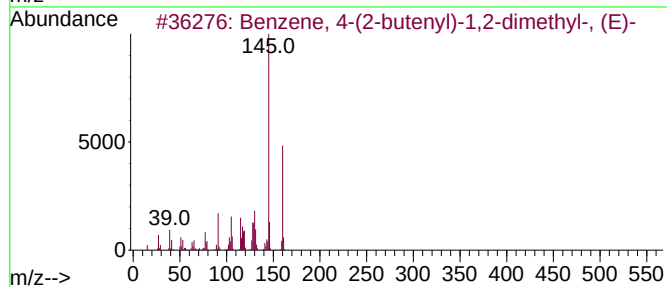
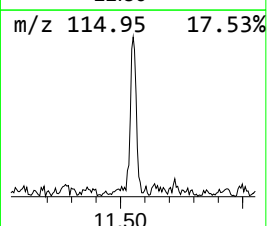
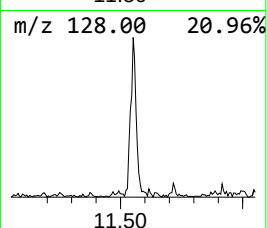
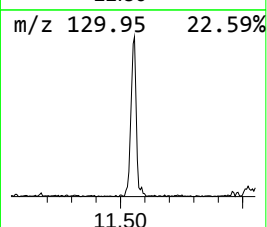
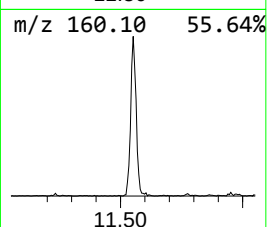
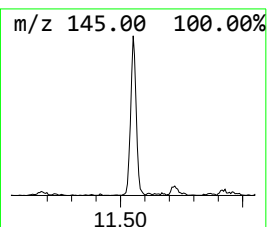
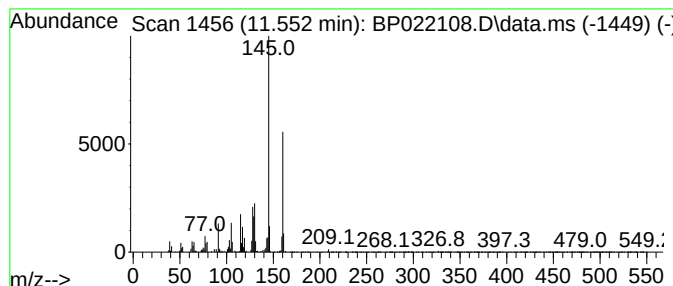
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.552	2.69 ng/ul	101500	Naphthalene-d8	10.458

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94
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-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	93
5			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022108.D
Acq On : 28 Sep 2024 17:07
Operator : RC/JU
Sample : P3910-08MEDL 20X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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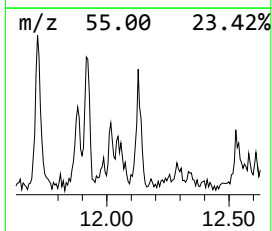
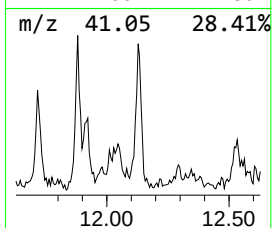
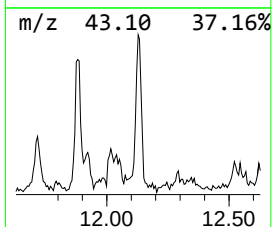
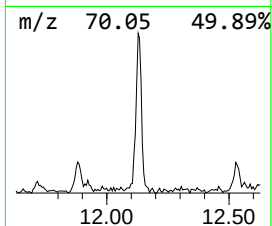
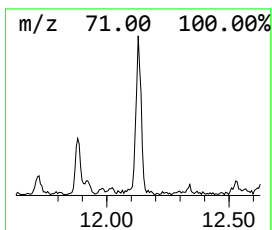
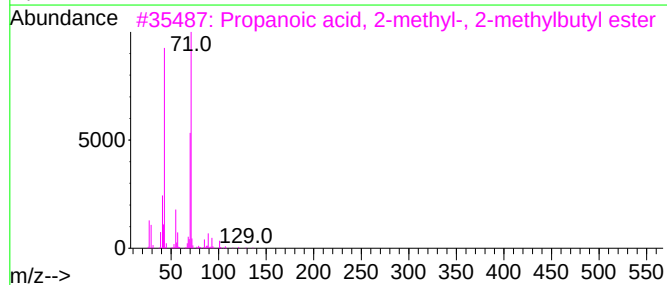
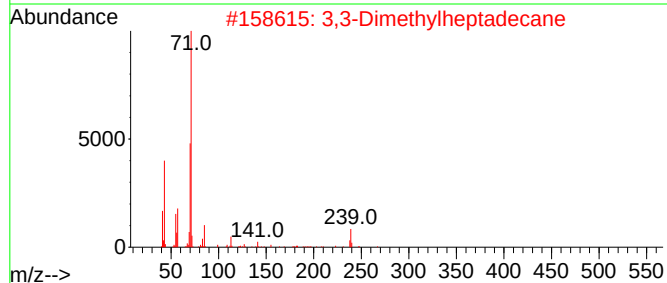
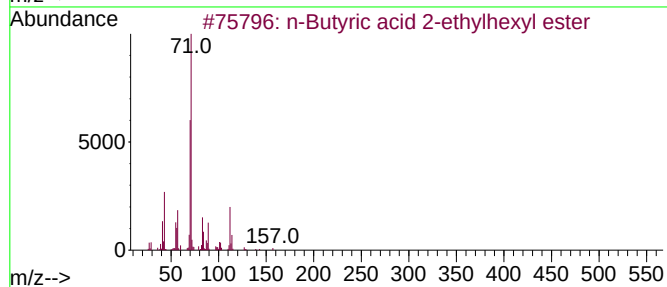
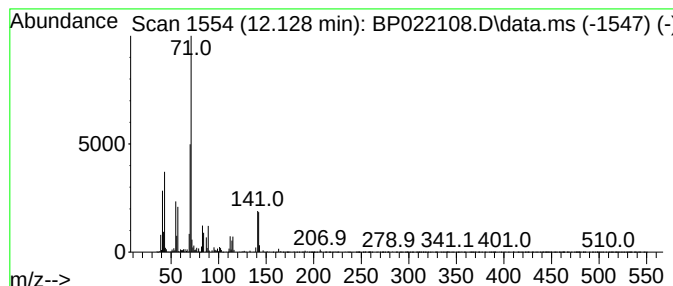
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Butyric acid 2-ethylhexyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.128	2.26 ng/ul	85452	Naphthalene-d8	10.458

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Butyric acid 2-ethylhexyl ester	200	C12H24O2	025415-84-3	72
2			3,3-Dimethylheptadecane	268	C19H40	1000360-43-3	53
3			Propanoic acid, 2-methyl-, 2-met...	158	C9H18O2	002445-69-4	53
4			Propanoic acid, 2-methyl-, 2-met...	158	C9H18O2	002445-69-4	53
5			2-Propyl-1-pentanol, 2-methylpro...	200	C12H24O2	1000447-36-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022108.D
Acq On : 28 Sep 2024 17:07
Operator : RC/JU
Sample : P3910-08MEDL 20X
Misc :
ALS Vial : 14 Sample Multiplier: 1

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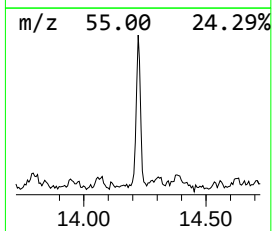
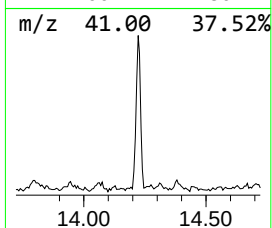
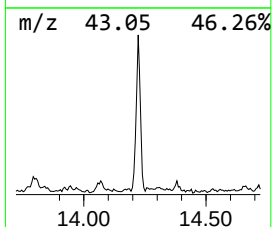
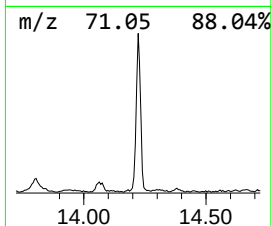
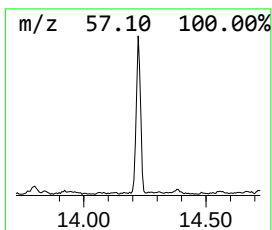
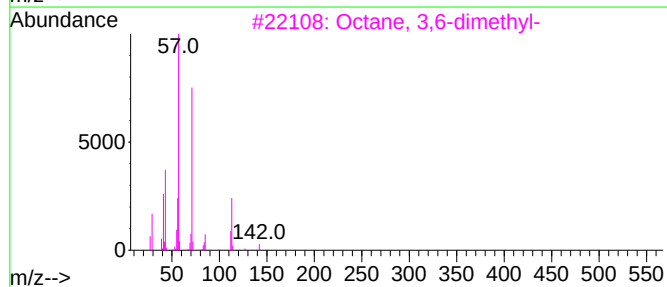
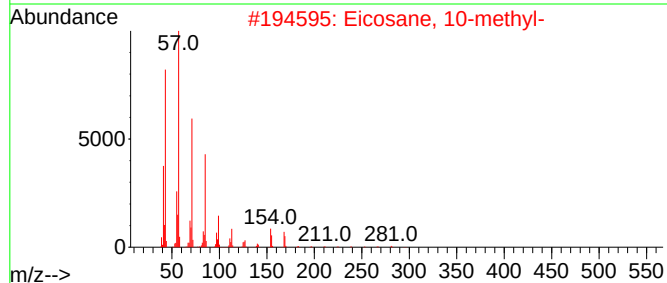
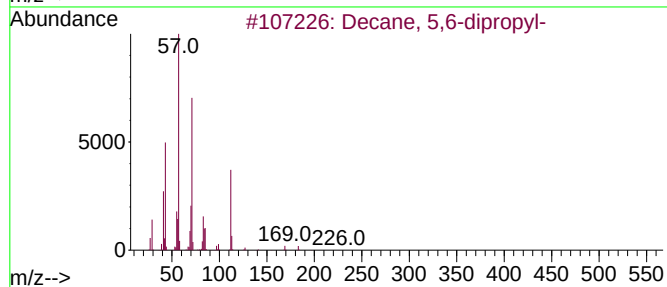
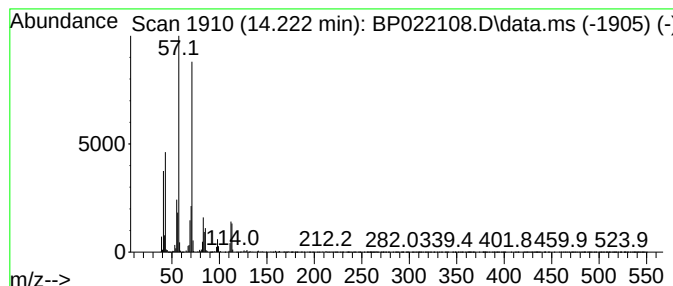
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Quant Title : SVOA CALIBRATION

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	4.09 ng/ul	213111	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	91
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	64
3			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
4			Propanal, 2-propenylhydrazone	112	C6H12N2	019031-78-8	58
5			Sulfurous acid, isobutyl 2-penty...	208	C9H20O3S	1000309-15-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022108.D
Acq On : 28 Sep 2024 17:07
Operator : RC/JU
Sample : P3910-08MEDL 20X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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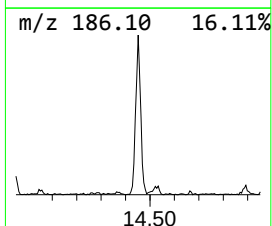
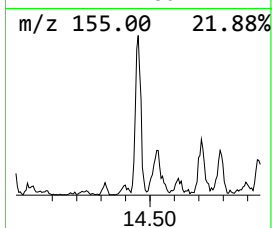
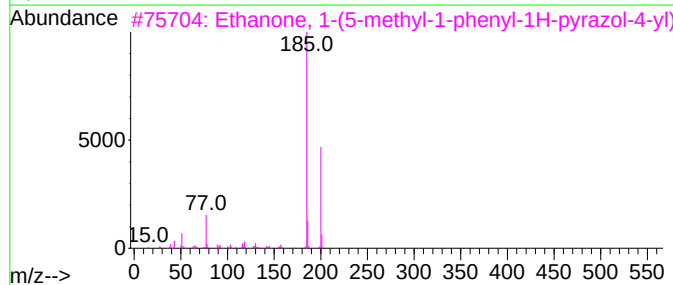
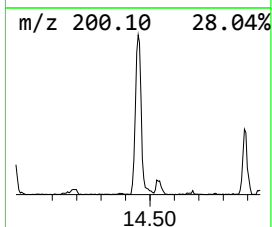
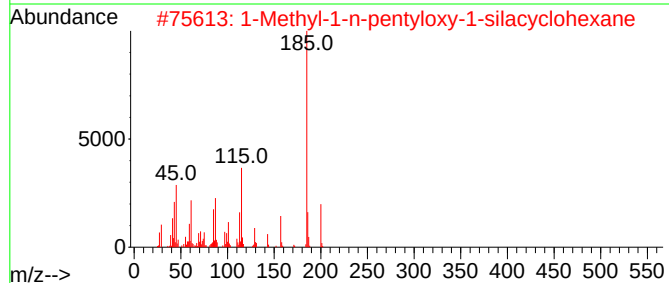
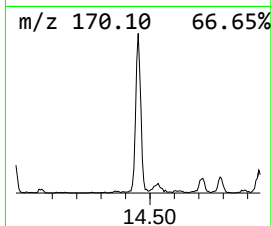
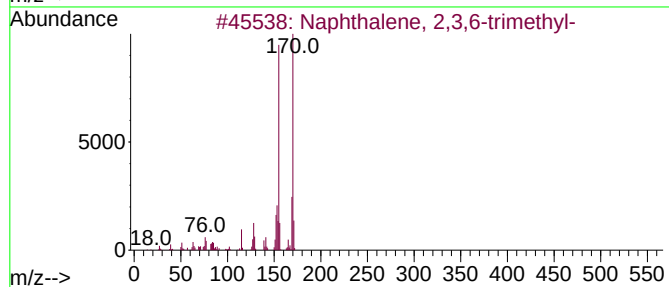
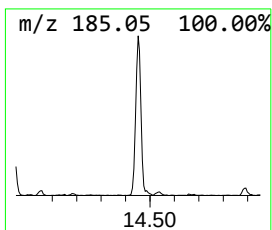
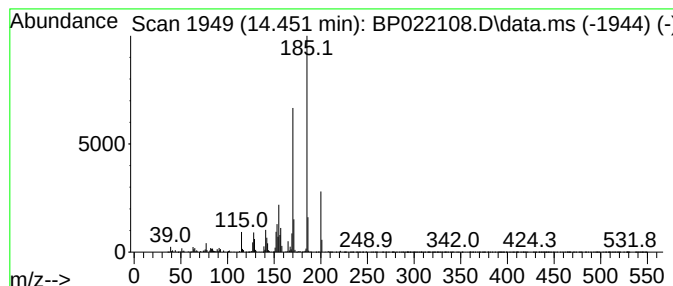
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Naphthalene, 2,3,6-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.451	3.04 ng/ul	158255	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	51
2			1-Methyl-1-n-pentyloxy-1-silacyc...	200	C11H24OSi	232270-61-0	43
3			Ethanone, 1-(5-methyl-1-phenyl-1...	200	C12H12N2O	006123-63-3	43
4			Phenol, 4-(phenylamino)-	185	C12H11NO	000122-37-2	38
5			2,3-Dihydro-8-methylfuro(2,3-b)q...	185	C12H11NO	064124-82-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022108.D
Acq On : 28 Sep 2024 17:07
Operator : RC/JU
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Misc :
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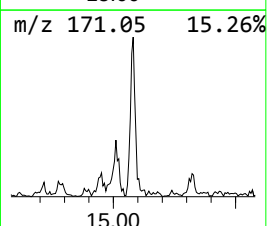
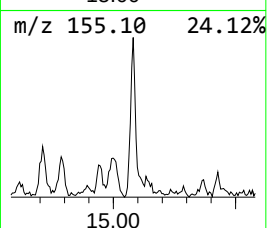
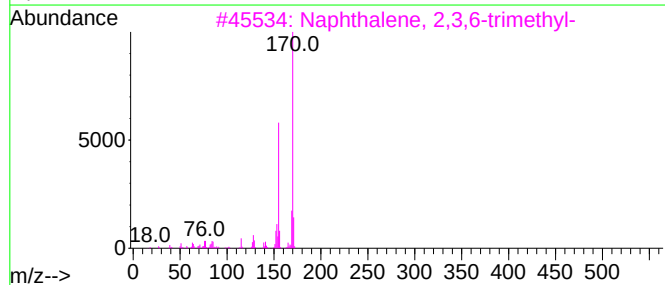
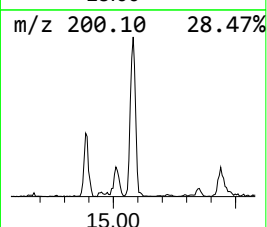
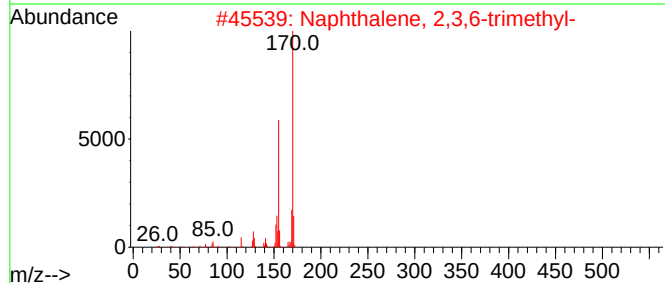
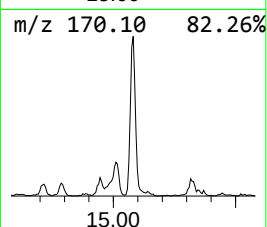
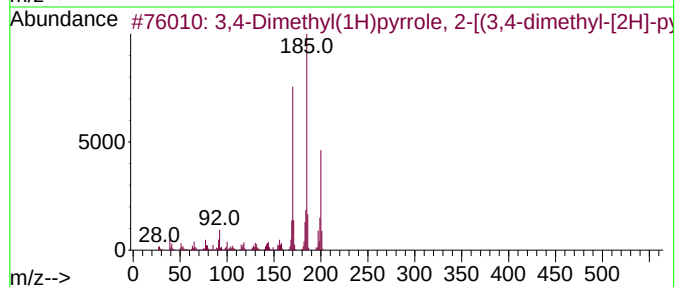
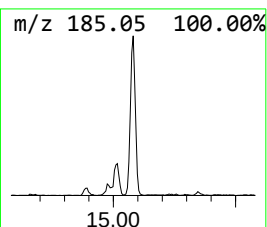
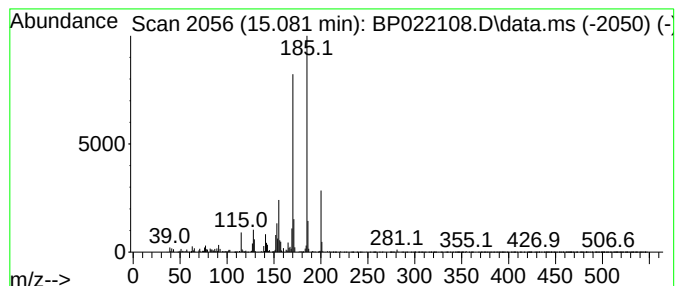
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 3,4-Dimethyl(1H)pyrrole, 2-... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.081	3.29 ng/ul	171387	Acenaphthene-d10	14.310

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,4-Dimethyl(1H)pyrrole, 2-[(3,4...	200	C13H16N2	065539-79-9	87
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	55
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	55
4			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	50
5			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022108.D
Acq On : 28 Sep 2024 17:07
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Sample : P3910-08MEDL 20X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	7.334	5.4	ng/ul	144383	1	7.699	535773	20.0
1-Hexanol, 2-et...	7.764	2.1	ng/ul	56575	1	7.699	535773	20.0
(DEL) Alkane: S...	8.911	3.8	ng/ul	101342	1	7.699	535773	20.0
Benzene, 4-(2-b...	11.552	2.7	ng/ul	101500	2	10.458	755558	20.0
n-Butyric acid ...	12.128	2.3	ng/ul	85452	2	10.458	755558	20.0
(DEL) Alkane: S...	14.222	4.1	ng/ul	213111	3	14.310	1040860	20.0
Naphthalene, 2,...	14.451	3.0	ng/ul	158255	3	14.310	1040860	20.0
3,4-Dimethyl(1H...	15.081	3.3	ng/ul	171387	3	14.310	1040860	20.0