

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018434.D
 Acq On : 28 Nov 2023 11:20
 Operator : MA/JU
 Sample : 05416-02
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AB3

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-BP112223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018434.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.269	27	31	38	rVB	86230	118218	1.67%	0.123%
2	3.528	71	75	84	rVB2	85240	154654	2.18%	0.161%
3	3.687	97	102	112	rBV	704775	1015923	14.35%	1.057%
4	4.022	155	159	167	rVB2	44815	69751	0.98%	0.073%
5	4.987	320	323	328	rVV	48434	69417	0.98%	0.072%
6	5.040	328	332	337	rVV	50792	79792	1.13%	0.083%
7	5.205	355	360	365	rVB	51949	72559	1.02%	0.075%
8	5.869	466	473	478	rVV2	109746	177653	2.51%	0.185%
9	6.128	508	517	524	rBV4	32798	74977	1.06%	0.078%
10	6.346	548	554	561	rVB5	35373	75531	1.07%	0.079%
11	6.493	570	579	582	rVV2	46649	102424	1.45%	0.107%
12	6.522	582	584	591	rVB2	55501	80240	1.13%	0.083%
13	6.757	616	624	628	rVV5	43930	99010	1.40%	0.103%
14	6.840	628	638	644	rVV7	48567	139047	1.96%	0.145%
15	6.904	646	649	653	rVV4	40560	75934	1.07%	0.079%
16	7.022	660	669	676	rVV	1275667	2102063	29.68%	2.187%
17	7.193	691	698	704	rBV	1139331	1739776	24.57%	1.810%
18	7.387	724	731	740	rVV	1316427	2103352	29.70%	2.188%
19	7.487	740	748	756	rVB4	127024	280808	3.97%	0.292%
20	7.857	803	811	819	rVV	1639771	2669124	37.69%	2.776%
21	8.522	916	924	926	rBV4	32479	66089	0.93%	0.069%
22	8.563	926	931	939	rVV	1595237	2497570	35.27%	2.598%
23	9.016	1002	1008	1016	rBV	1236493	1922044	27.14%	1.999%
24	9.093	1016	1021	1031	rVB	193798	306328	4.33%	0.319%
25	9.451	1072	1082	1087	rVV9	27748	84554	1.19%	0.088%
26	9.740	1121	1131	1137	rBV	830582	1397864	19.74%	1.454%
27	10.093	1182	1191	1195	rBV5	41255	83274	1.18%	0.087%
28	10.275	1213	1222	1231	rVB	1609616	2672692	37.74%	2.780%
29	10.651	1277	1286	1292	rVV	2199899	4022435	56.80%	4.184%
30	10.704	1292	1295	1301	rVV	151673	236593	3.34%	0.246%
31	10.793	1302	1310	1323	rVV	1389332	2454789	34.66%	2.553%
32	11.581	1436	1444	1452	rVV4	41716	83926	1.19%	0.087%
33	11.692	1457	1463	1469	rBV	233737	353825	5.00%	0.368%
34	12.087	1524	1530	1537	rBV	312472	454229	6.41%	0.472%
35	12.239	1549	1556	1562	rVV3	37073	84322	1.19%	0.088%
36	12.316	1562	1569	1577	rVB	313235	501495	7.08%	0.522%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
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37	12.534	1600	1606	1612	rVV	244611	369955	5.22%	0.385%
38	12.775	1641	1647	1652	rVB5	37862	73505	1.04%	0.076%
39	12.898	1664	1668	1672	rVV	52976	83742	1.18%	0.087%
40	13.039	1688	1692	1697	rVV	156396	215841	3.05%	0.225%
41	13.328	1734	1741	1746	rBV	455518	622069	8.78%	0.647%
42	13.398	1748	1753	1760	rVB	82118	141354	2.00%	0.147%
43	13.651	1791	1796	1797	rBV	96353	117464	1.66%	0.122%
44	13.810	1818	1823	1828	rVV	179437	309553	4.37%	0.322%
45	13.869	1828	1833	1835	rVV	165283	262348	3.70%	0.273%
46	13.904	1835	1839	1846	rVV	2825180	3915812	55.29%	4.073%
47	13.986	1848	1853	1857	rVV	258751	374092	5.28%	0.389%
48	14.051	1857	1864	1867	rVV2	106663	219689	3.10%	0.229%
49	14.186	1876	1887	1898	rBV	3344990	5093946	71.93%	5.299%
50	14.398	1919	1923	1929	rVB	402464	501436	7.08%	0.522%
51	14.492	1929	1939	1945	rBV	2989477	4748911	67.06%	4.940%
52	14.581	1950	1954	1958	rVB	46992	67589	0.95%	0.070%
53	14.681	1965	1971	1981	rBV	654461	1070296	15.11%	1.113%
54	14.892	2002	2007	2014	rVV	193858	346890	4.90%	0.361%
55	14.963	2014	2019	2023	rVV	62440	95596	1.35%	0.099%
56	15.022	2025	2029	2036	rVB5	65538	109892	1.55%	0.114%
57	15.110	2036	2044	2048	rBV3	55812	124996	1.77%	0.130%
58	15.163	2048	2053	2057	rVB	70407	98136	1.39%	0.102%
59	15.281	2067	2073	2077	rBV3	109988	172013	2.43%	0.179%
60	15.351	2081	2085	2090	rVB	372040	445792	6.29%	0.464%
61	15.480	2101	2107	2113	rVV	3998068	5975308	84.37%	6.215%
62	15.528	2113	2115	2120	rVB	84032	98133	1.39%	0.102%
63	15.598	2120	2127	2135	rBV	528840	785411	11.09%	0.817%
64	15.757	2150	2154	2159	rVV	133788	190371	2.69%	0.198%
65	15.833	2159	2167	2174	rVB	201839	321944	4.55%	0.335%
66	15.980	2187	2192	2200	rVB	166550	338981	4.79%	0.353%
67	16.069	2200	2207	2213	rVB2	73609	128347	1.81%	0.134%
68	16.222	2227	2233	2234	rBV	391191	513531	7.25%	0.534%
69	16.239	2234	2236	2242	rVB	437167	556989	7.87%	0.579%
70	16.780	2323	2328	2332	rBV3	74517	130566	1.84%	0.136%
71	16.892	2343	2347	2355	rVB4	71679	139282	1.97%	0.145%
72	16.975	2356	2361	2364	rBV2	75357	130527	1.84%	0.136%
73	17.016	2364	2368	2373	rVV	370872	473545	6.69%	0.493%
74	17.069	2373	2377	2384	rVB4	79443	129645	1.83%	0.135%
75	17.245	2400	2407	2411	rBV	3010490	4387559	61.95%	4.564%
76	17.286	2411	2414	2418	rVV	435936	562870	7.95%	0.585%

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77	17.345	2418	2424	2434	rVB	3668157	5401430	76.27%	5.619%
78	17.427	2434	2438	2444	rBV6	38857	69010	0.97%	0.072%
79	17.557	2456	2460	2467	rVB3	77527	118188	1.67%	0.123%
80	17.763	2490	2495	2503	rVV	373399	476680	6.73%	0.496%
81	18.116	2551	2555	2557	rBV	89103	111104	1.57%	0.116%
82	18.151	2557	2561	2568	rVB2	402088	680645	9.61%	0.708%
83	18.292	2581	2585	2589	rBV	114180	163994	2.32%	0.171%
84	18.339	2589	2593	2599	rVB	138611	186353	2.63%	0.194%
85	18.439	2605	2610	2620	rBV	335084	472841	6.68%	0.492%
86	18.604	2635	2638	2643	rVB	99862	122485	1.73%	0.127%
87	19.010	2703	2707	2712	rBV	77511	114684	1.62%	0.119%
88	19.063	2712	2716	2719	rVV	291676	336237	4.75%	0.350%
89	19.104	2719	2723	2727	rVB	55914	73031	1.03%	0.076%
90	19.169	2731	2734	2739	rVB2	70436	97154	1.37%	0.101%
91	19.274	2748	2752	2756	rVB2	68908	106388	1.50%	0.111%
92	19.398	2770	2773	2778	rVB3	101596	121558	1.72%	0.126%
93	19.598	2801	2807	2811	rBV	4431660	6042605	85.33%	6.285%
94	19.633	2811	2813	2818	rVB	226923	261628	3.69%	0.272%
95	20.174	2902	2905	2911	rBV	218128	259962	3.67%	0.270%
96	20.692	2990	2993	3002	rVB	191034	262198	3.70%	0.273%
97	21.198	3073	3079	3089	rBV2	324457	685569	9.68%	0.713%
98	21.404	3108	3114	3123	rBV	3252789	4733412	66.84%	4.924%
99	23.809	3514	3523	3540	rVB2	2431949	7081868	100.00%	7.367%
100	23.986	3544	3553	3566	rVB	1739581	5190754	73.30%	5.399%

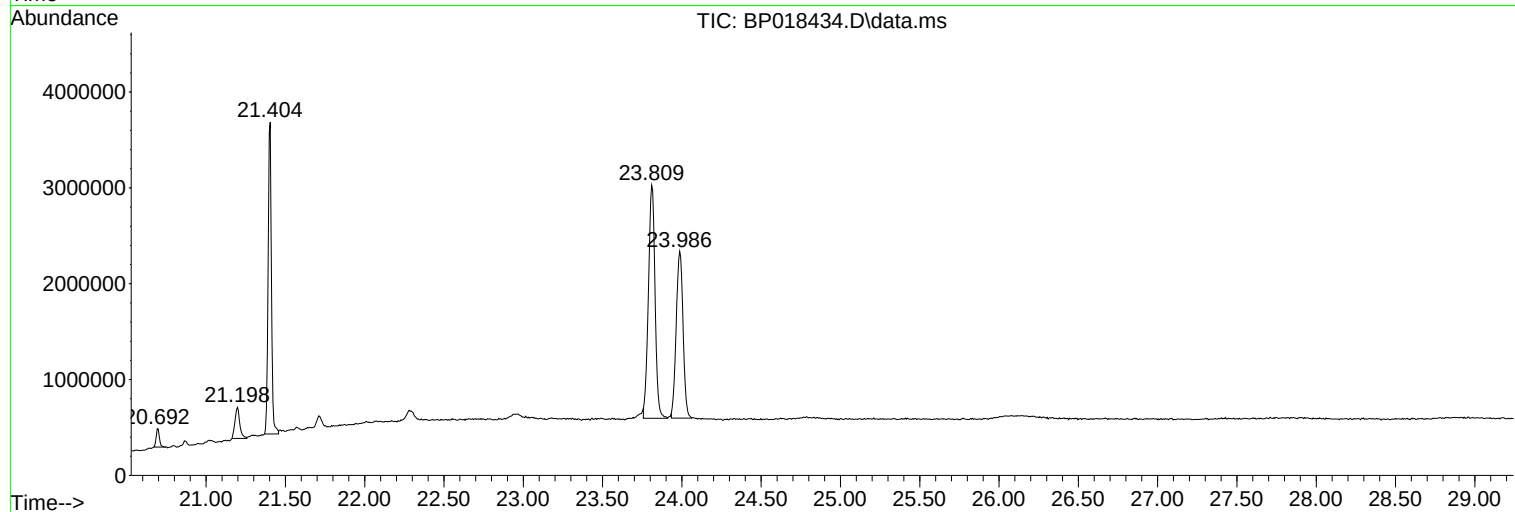
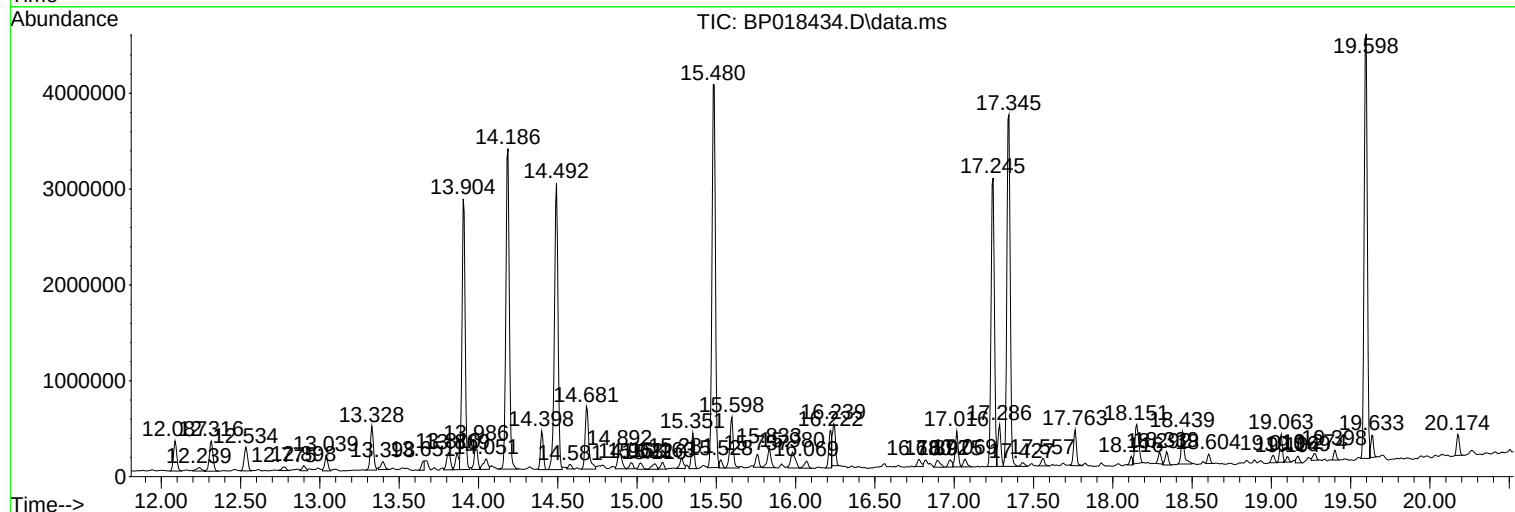
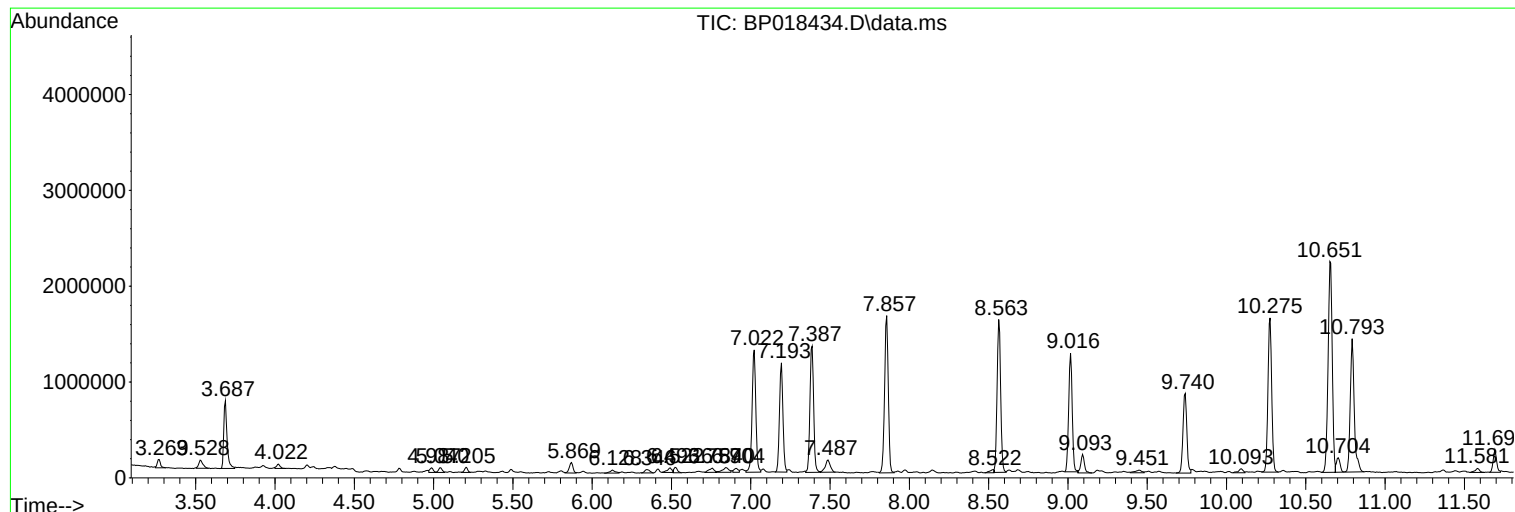
Sum of corrected areas: 96135986

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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



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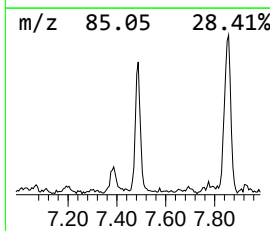
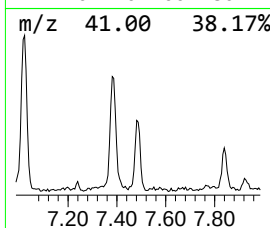
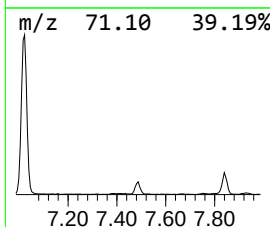
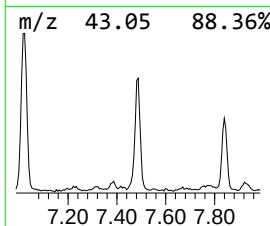
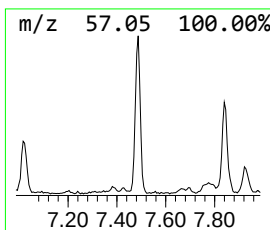
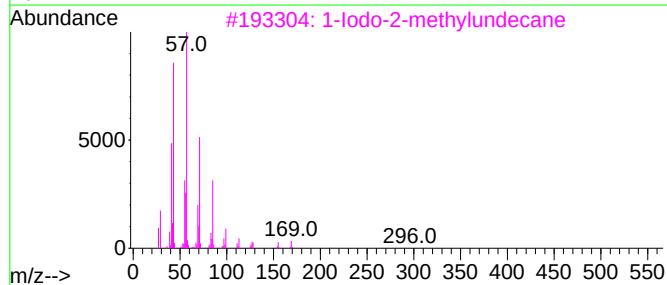
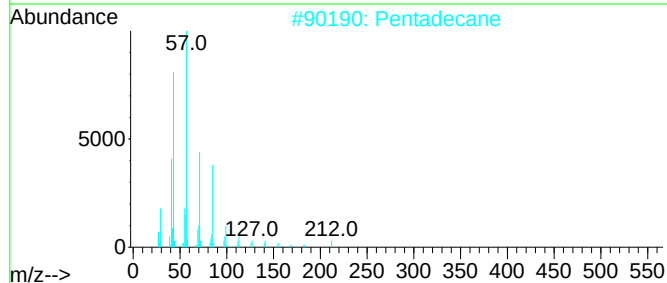
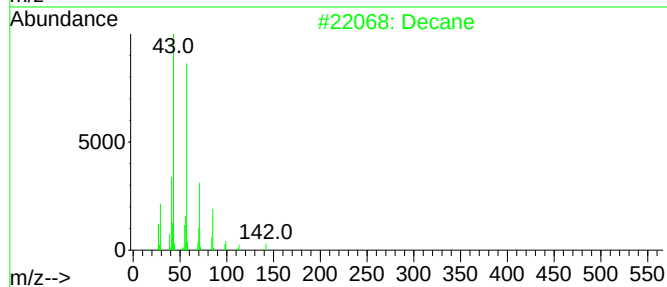
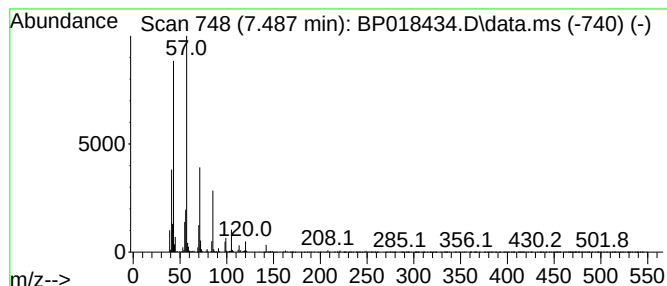
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.487	2.10 ng/ul	280808	1,4-Dichlorobenzene-d4	7.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	81
2			Pentadecane	212	C15H32	000629-62-9	72
3			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	72
4			Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	72
5			Hexadecane	226	C16H34	000544-76-3	64



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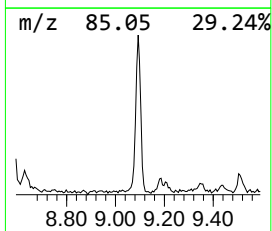
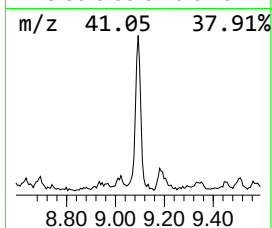
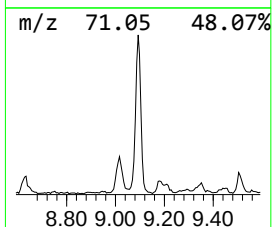
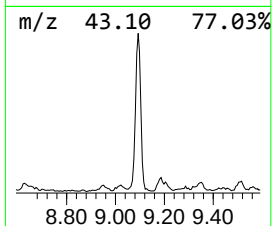
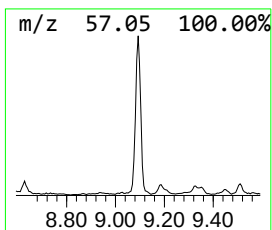
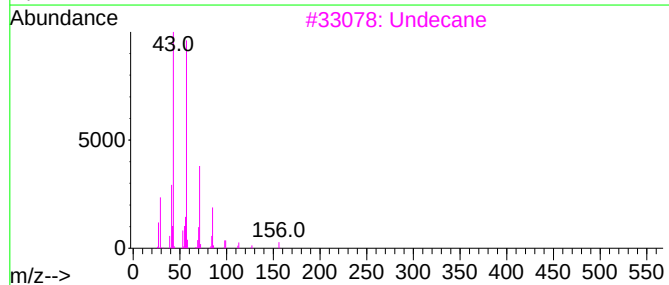
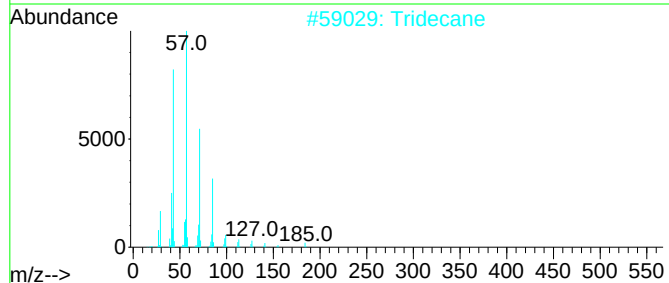
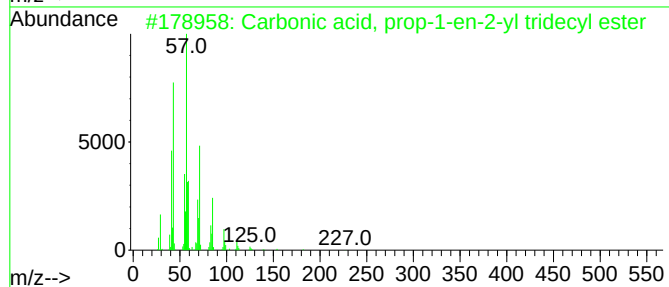
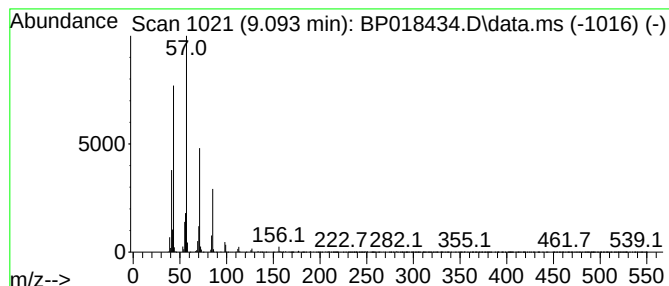
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Carbonic acid, prop-1-en-2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.093	2.30 ng/ul	306328	1,4-Dichlorobenzene-d4	7.857

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	78
2			Tridecane	184	C13H28	000629-50-5	78
3			Undecane	156	C11H24	001120-21-4	76
4			Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	72
5			Undecane	156	C11H24	001120-21-4	58



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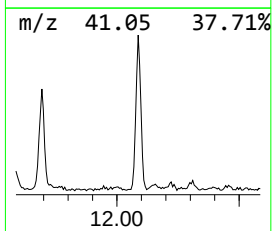
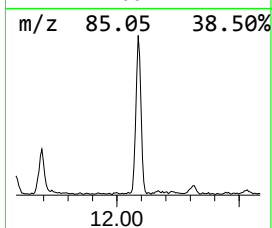
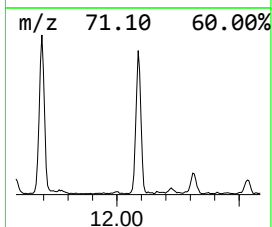
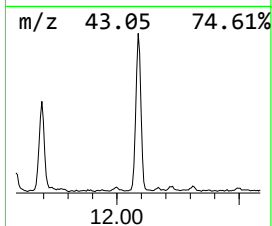
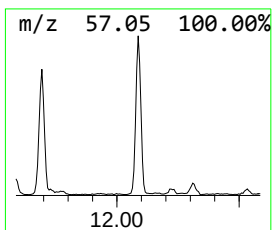
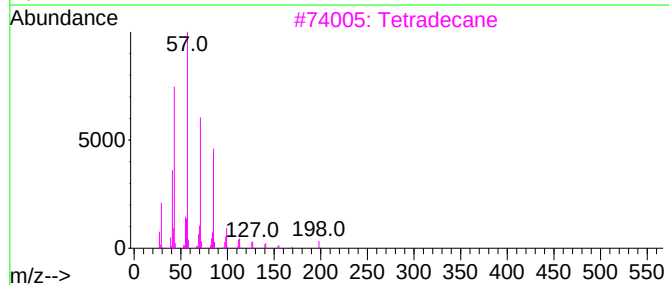
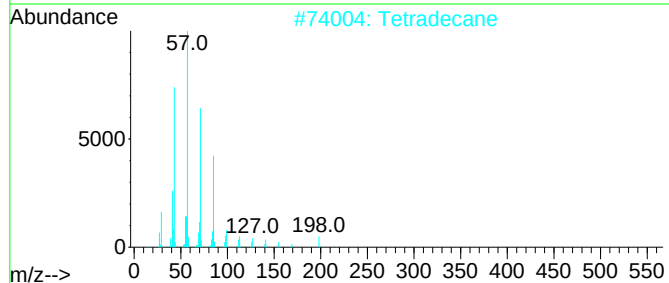
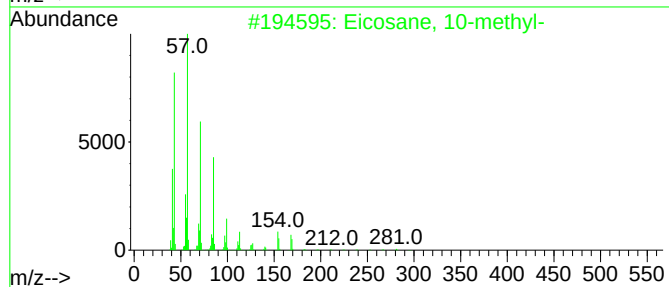
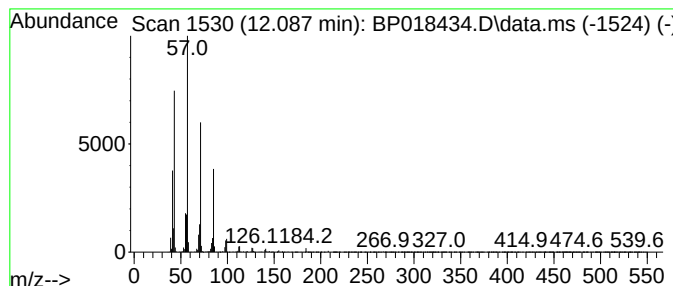
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.087	2.26 ng/ul	454229	Naphthalene-d8	10.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
2		Tetradecane	198	C14H30	000629-59-4	90
3		Tetradecane	198	C14H30	000629-59-4	90
4		Tridecane	184	C13H28	000629-50-5	81
5		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
Operator : MA/JU
Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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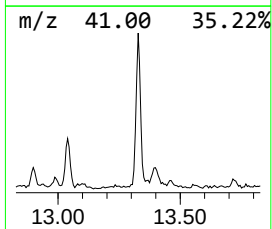
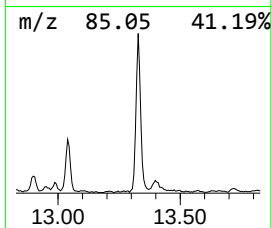
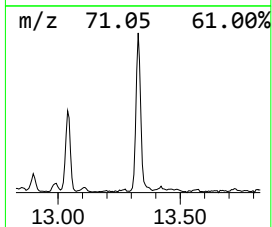
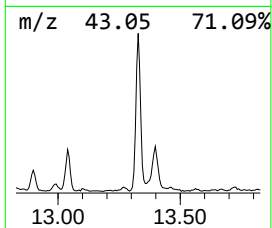
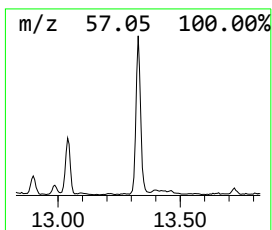
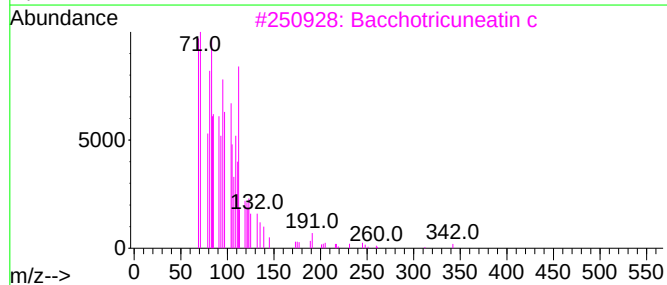
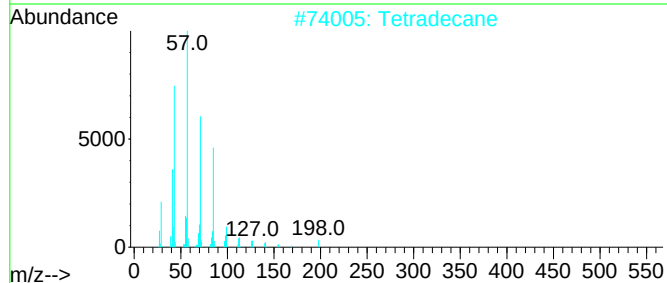
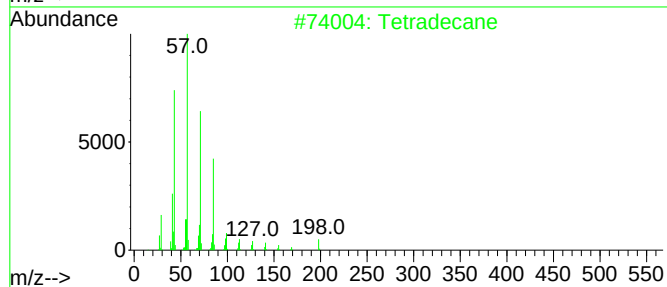
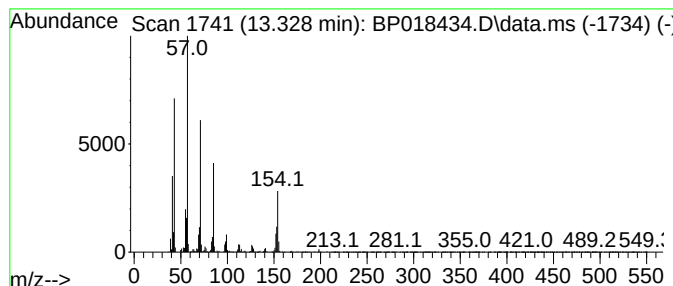
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.328	2.62 ng/ul	622069	Acenaphthene-d10	14.492

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	89
2			Tetradecane	198	C14H30	000629-59-4	62
3			Bacchotricuneatin c	342	C20H22O5	066563-30-2	60
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	58
5			Tetradecane	198	C14H30	000629-59-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
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Acq On : 28 Nov 2023 11:20
Operator : MA/JU
Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

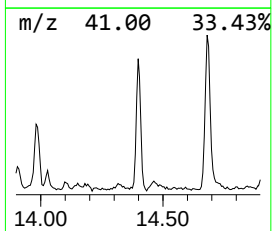
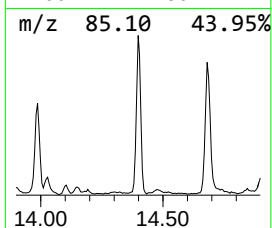
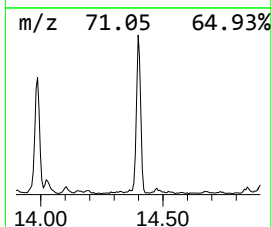
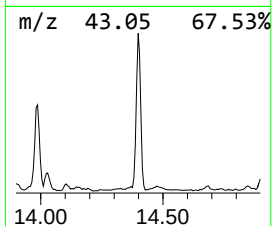
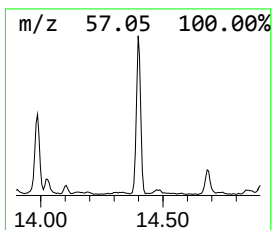
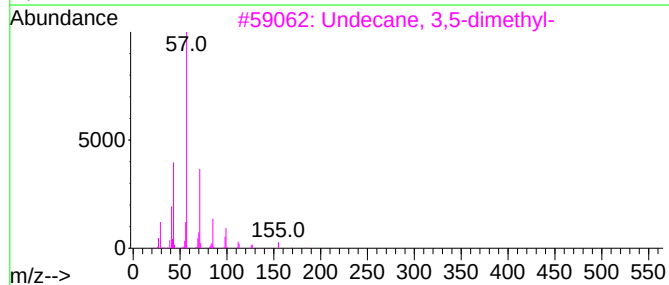
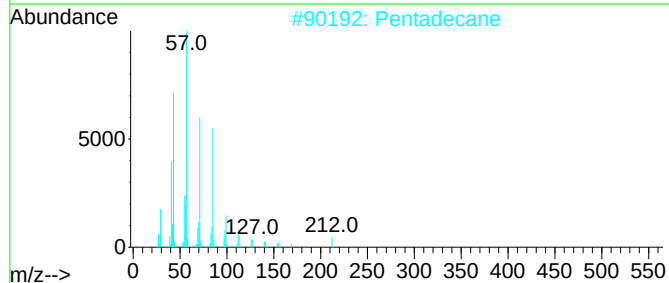
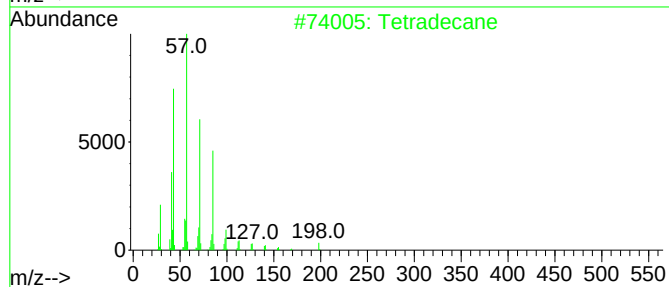
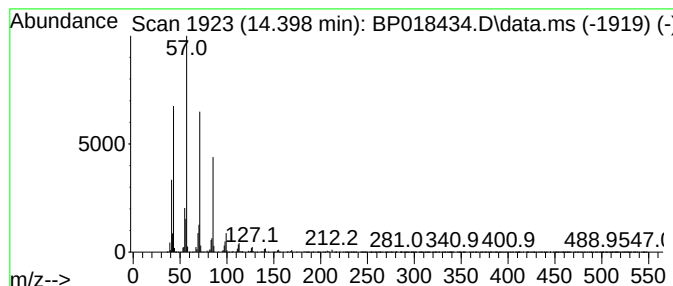
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Quant Title : SVOA CALIBRATION

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.	
14.398	2.11 ng/ul	501436	Acenaphthene-d10	14.492	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane	198	C14H30	000629-59-4	91
2	Pentadecane	212	C15H32	000629-62-9	90
3	Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	87
4	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	86
5	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
Operator : MA/JU
Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

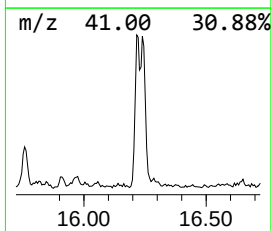
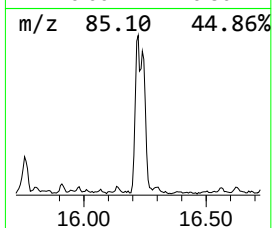
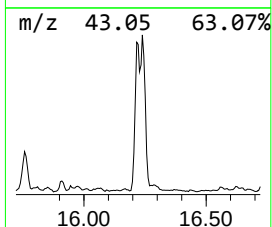
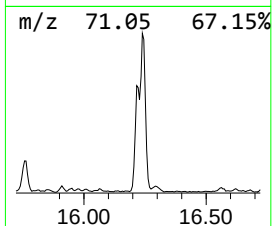
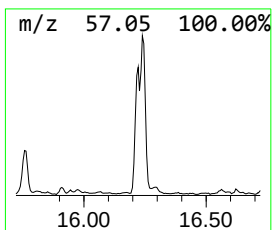
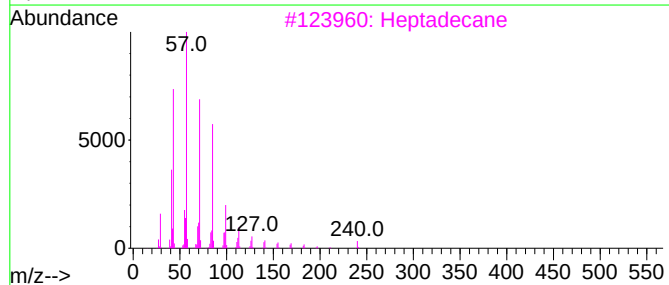
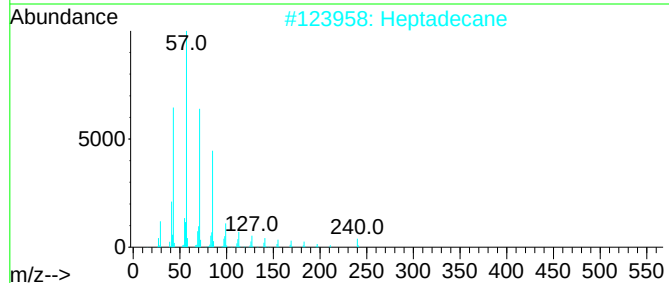
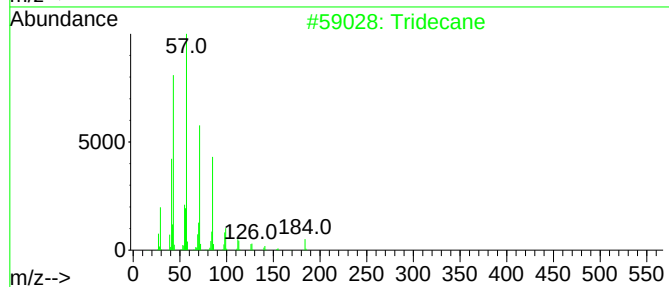
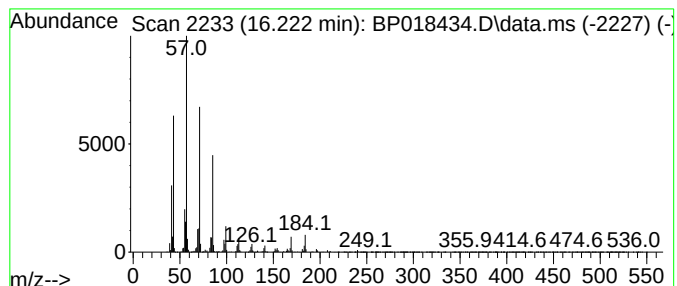
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.222	2.34 ng/ul	513531	Phenanthrene-d10		17.245	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	94
2	Heptadecane		240	C17H36	000629-78-7	93
3	Heptadecane		240	C17H36	000629-78-7	90
4	Tridecane, 7-propyl-		226	C16H34	055045-09-5	90
5	Eicosane		282	C20H42	000112-95-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
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Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
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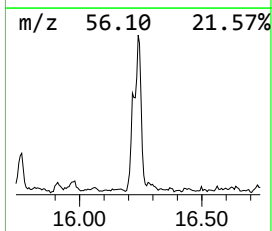
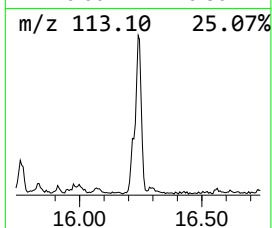
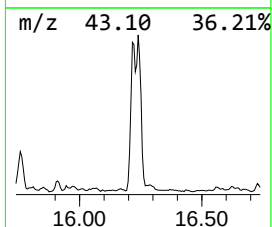
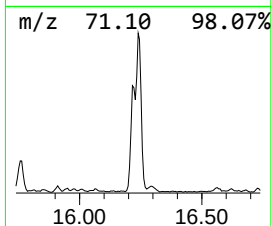
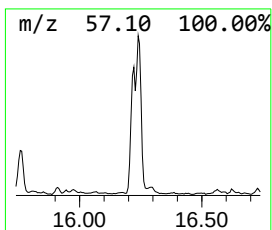
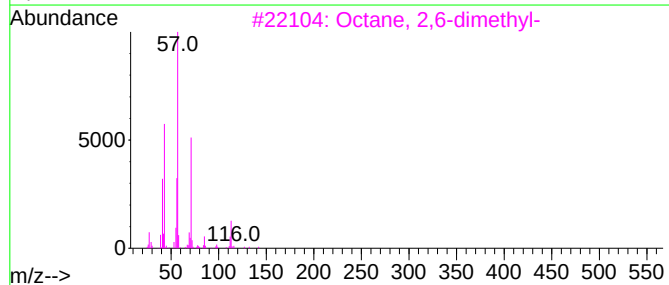
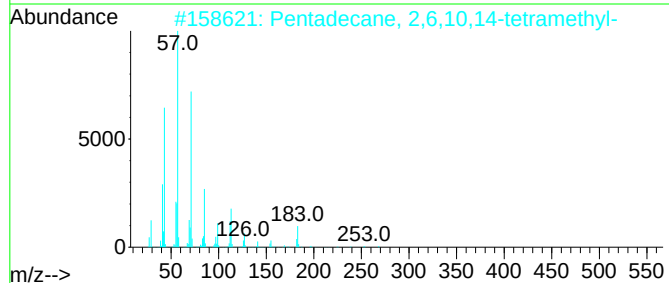
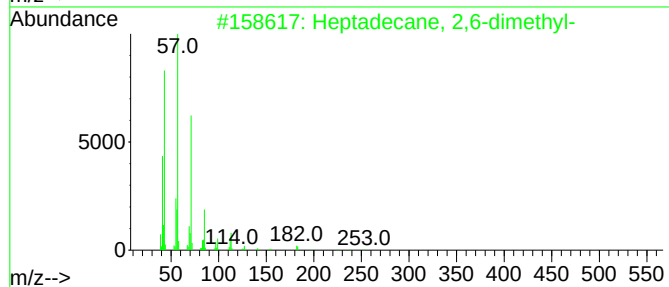
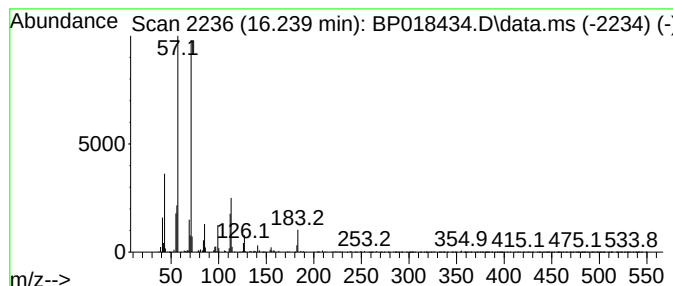
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.239	2.54 ng/ul	556989	Phenanthrene-d10	17.245

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	86
2		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	86
3		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	80
4		Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	78
5		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
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Sample : 05416-02
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Instrument :
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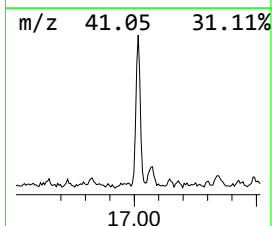
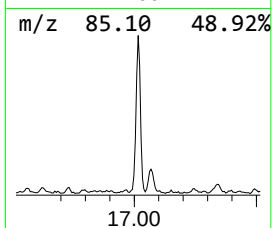
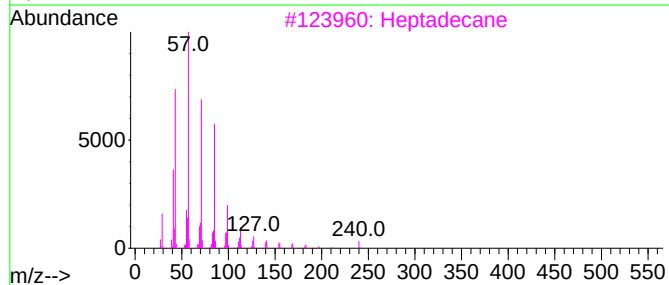
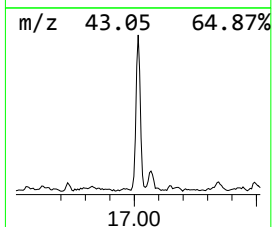
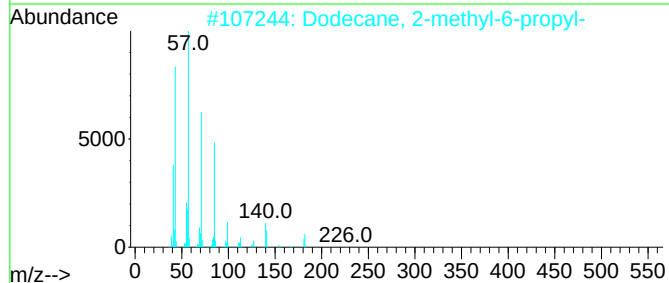
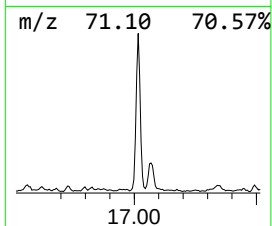
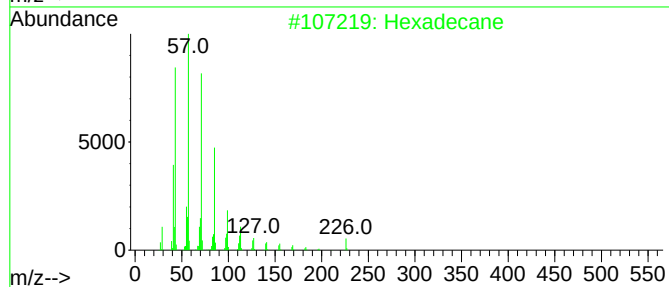
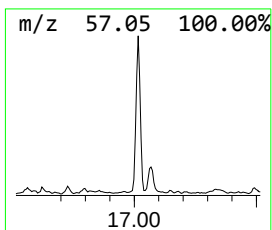
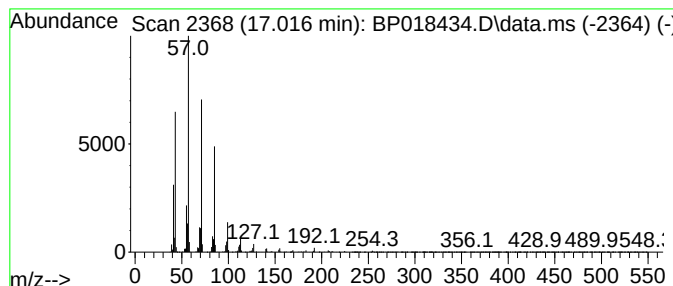
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.016	2.16 ng/ul	473545	Phenanthrene-d10	17.245

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	95
2		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
3		Heptadecane	240	C17H36	000629-78-7	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Pentadecane	212	C15H32	000629-62-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
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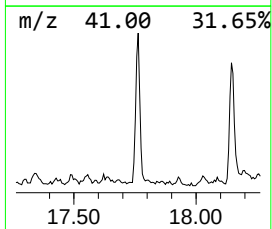
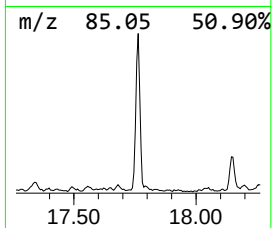
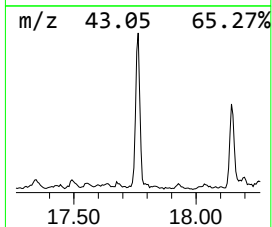
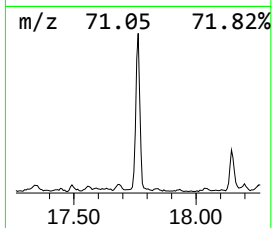
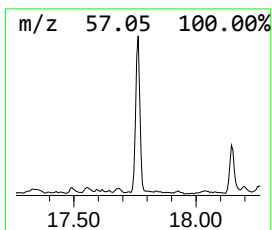
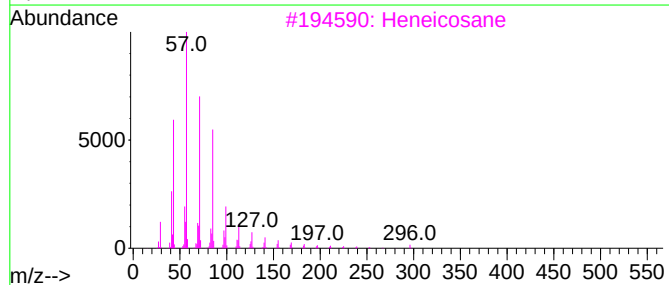
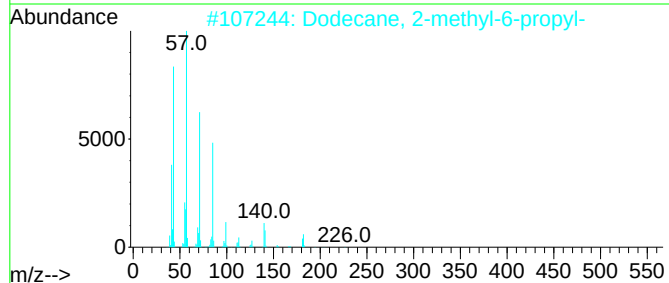
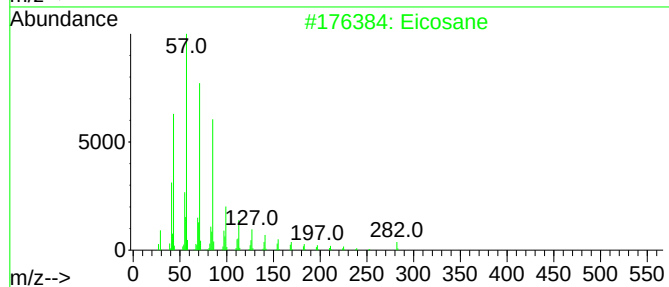
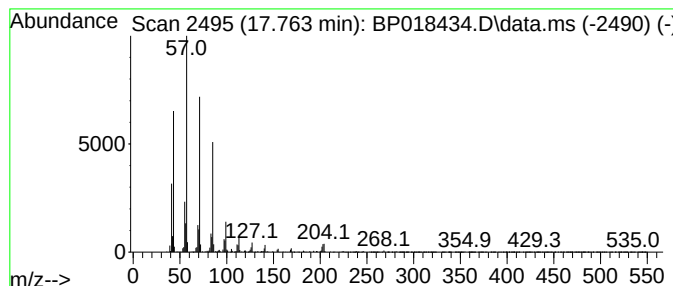
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.763	2.17 ng/ul	476680	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	95
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
3			Heneicosane	296	C21H44	000629-94-7	87
4			Octadecane	254	C18H38	000593-45-3	87
5			Tetradecane	198	C14H30	000629-59-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
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Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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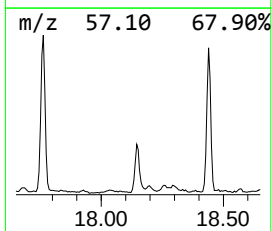
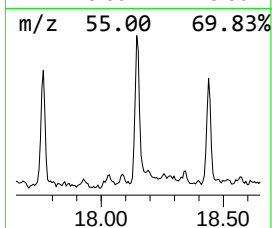
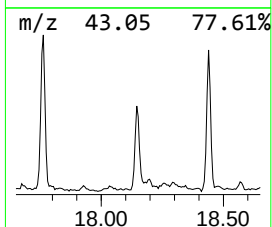
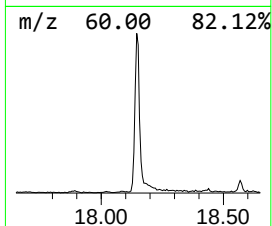
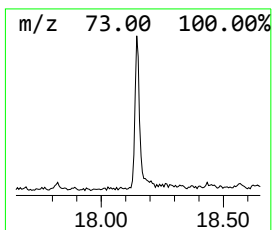
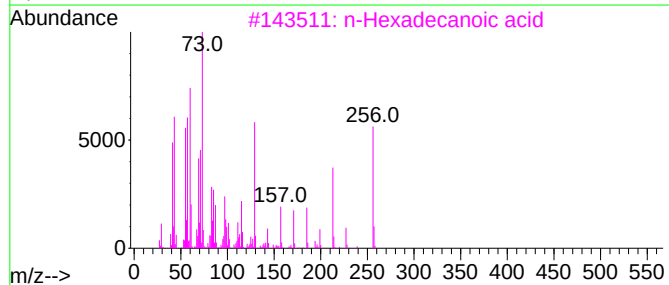
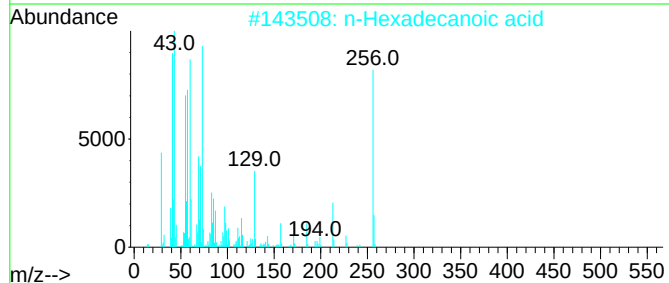
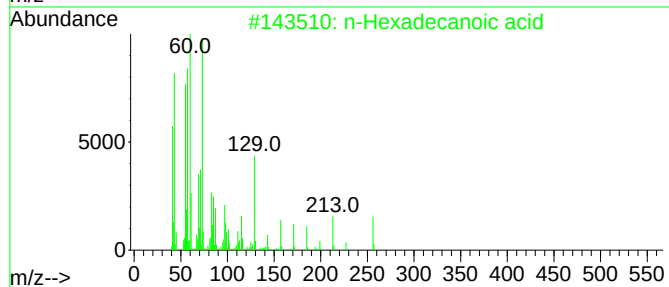
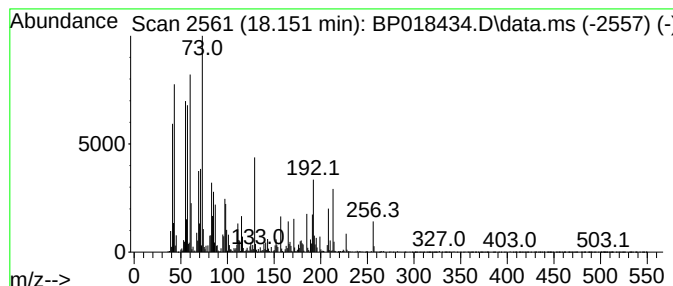
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.151	3.10 ng/ul	680645	Phenanthrene-d10	17.245

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5		Tridecanoic acid	214	C13H26O2	000638-53-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
Operator : MA/JU
Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB3

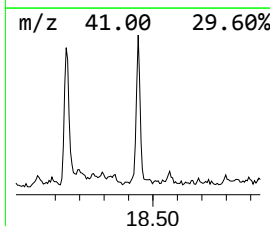
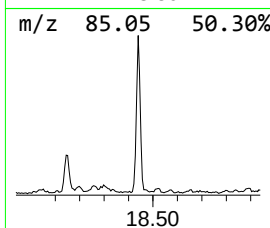
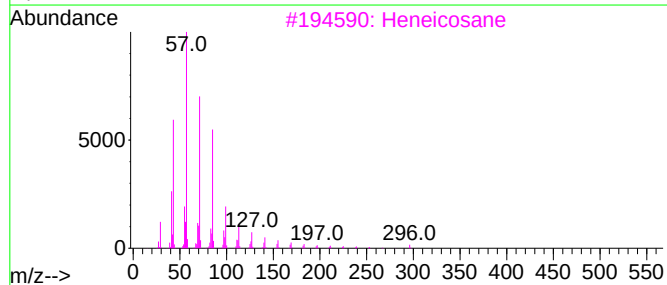
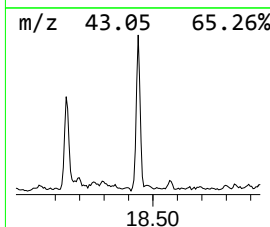
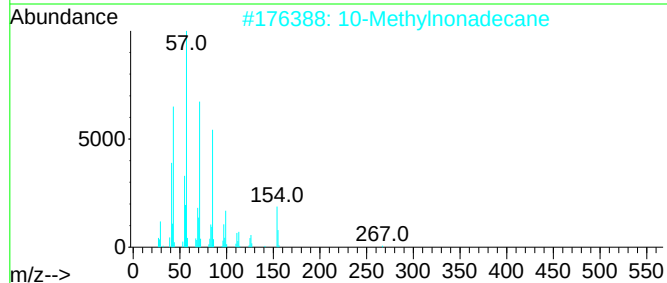
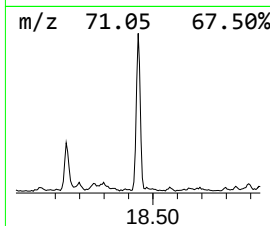
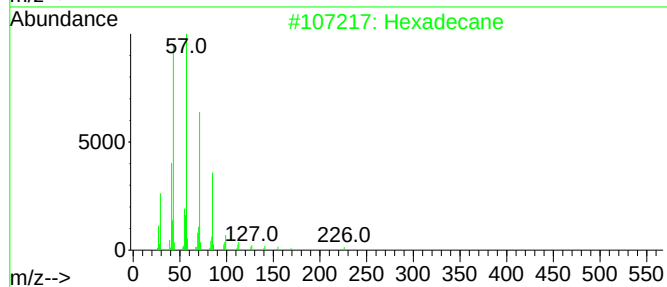
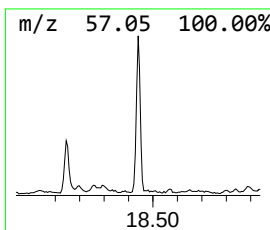
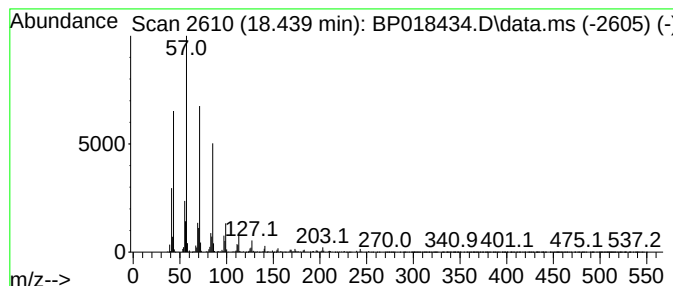
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.439	2.16 ng/ul	472841	Phenanthrene-d10	17.245

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	95
2		10-Methylnonadecane	282	C20H42	056862-62-5	91
3		Heneicosane	296	C21H44	000629-94-7	91
4		Heptadecane	240	C17H36	000629-78-7	91
5		Tridecane, 6-propyl-	226	C16H34	055045-10-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
Operator : MA/JU
Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB3

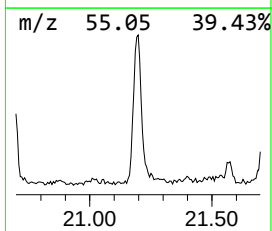
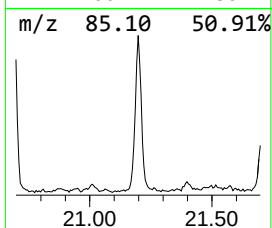
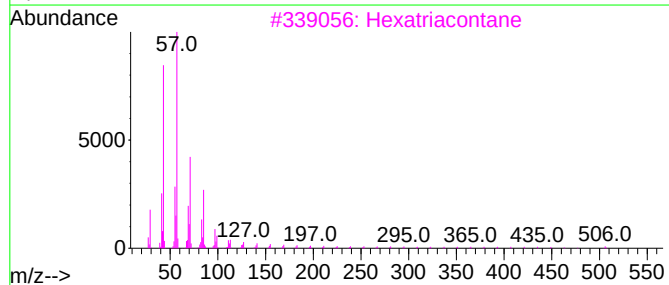
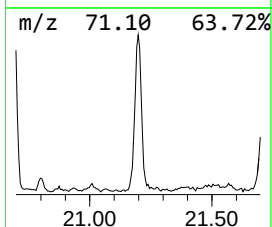
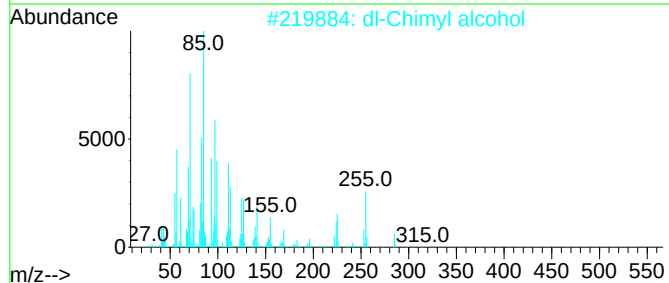
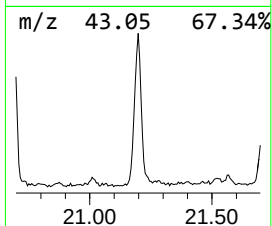
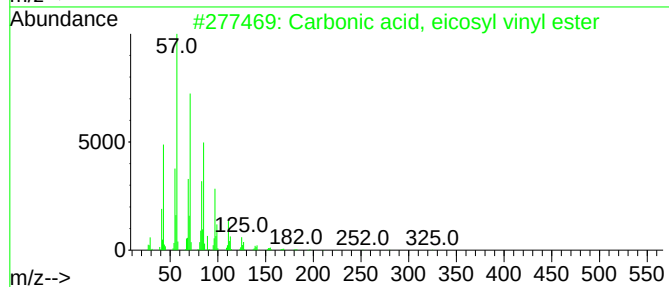
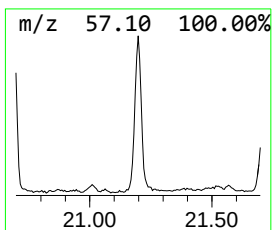
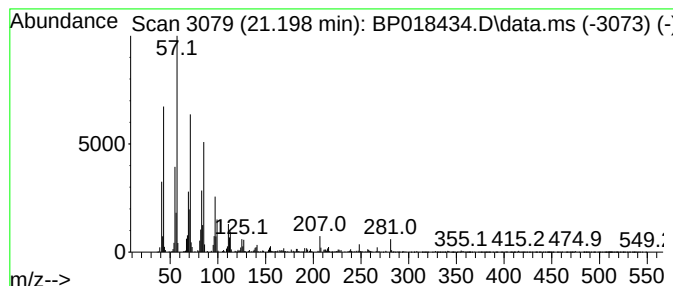
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Carbonic acid, eicosyl vinyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.198	2.90 ng/ul	685569	Chrysene-d12	21.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	94
2			dl-Chimyl alcohol	316	C19H40O3	006145-69-3	93
3			Hexatriacontane	507	C36H74	000630-06-8	91
4			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	91
5			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018434.D
Acq On : 28 Nov 2023 11:20
Operator : MA/JU
Sample : 05416-02
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.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...	7.487	2.1	ng/ul	280808	1	7.857	2669120	20.0
Carbonic acid, ...	9.093	2.3	ng/ul	306328	1	7.857	2669120	20.0
(DEL) Alkane: S...	12.087	2.3	ng/ul	454229	2	10.657	4022440	20.0
(DEL) Alkane: S...	13.328	2.6	ng/ul	622069	3	14.492	4748910	20.0
(DEL) Alkane: S...	14.398	2.1	ng/ul	501436	3	14.492	4748910	20.0
(DEL) Alkane: S...	16.222	2.3	ng/ul	513531	4	17.245	4387560	20.0
(DEL) Alkane: S...	16.239	2.5	ng/ul	556989	4	17.245	4387560	20.0
(DEL) Alkane: S...	17.016	2.2	ng/ul	473545	4	17.245	4387560	20.0
(DEL) Alkane: S...	17.763	2.2	ng/ul	476680	4	17.245	4387560	20.0
n-Hexadecanoic ...	18.151	3.1	ng/ul	680645	4	17.245	4387560	20.0
(DEL) Alkane: S...	18.439	2.2	ng/ul	472841	4	17.245	4387560	20.0
Carbonic acid, ...	21.198	2.9	ng/ul	685569	5	21.404	4733410	20.0