

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
 Data File : BP015998.D
 Acq On : 05 Jul 2023 04:07
 Operator : MA/JU
 Sample : 03244-22
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGG1

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015998.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.370	27	31	36	rVB	65312	83387	1.16%	0.104%
2	3.811	103	106	109	rBV3	10270	13644	0.19%	0.017%
3	3.834	109	110	120	rVV	93901	170782	2.37%	0.213%
4	3.917	121	124	130	rVB	45553	63655	0.88%	0.079%
5	4.028	139	143	150	rVB	84734	130411	1.81%	0.163%
6	4.134	158	161	165	rVB	25507	31939	0.44%	0.040%
7	4.375	198	202	209	rVB	64719	87952	1.22%	0.110%
8	4.799	270	274	281	rVB3	20359	34345	0.48%	0.043%
9	5.081	319	322	330	rVB2	15982	24418	0.34%	0.030%
10	5.999	473	478	484	rBV10	10700	22198	0.31%	0.028%
11	6.675	586	593	599	rVB3	20943	45622	0.63%	0.057%
12	7.116	665	668	671	rBV5	7250	12090	0.17%	0.015%
13	7.234	679	688	705	rBV	566042	1575658	21.87%	1.967%
14	7.364	705	710	721	rVV	700123	1308352	18.16%	1.634%
15	7.440	721	723	729	rVB	45211	67877	0.94%	0.085%
16	7.575	739	746	759	rBV	873618	1663967	23.10%	2.078%
17	8.034	818	824	839	rBV	1190230	2293721	31.84%	2.864%
18	8.240	853	859	865	rBV	57592	88848	1.23%	0.111%
19	8.328	870	874	880	rVB8	12207	21437	0.30%	0.027%
20	8.540	902	910	911	rBV8	9134	14971	0.21%	0.019%
21	8.793	945	953	967	rBV	593335	1716865	23.83%	2.144%
22	8.899	967	971	983	rVB	128757	285539	3.96%	0.357%
23	9.234	1020	1028	1045	rBV	739056	1624324	22.55%	2.028%
24	9.375	1047	1052	1065	rVB6	40658	99777	1.39%	0.125%
25	9.963	1145	1152	1171	rBV	442918	1271556	17.65%	1.588%
26	10.269	1196	1204	1206	rBV9	15935	34029	0.47%	0.042%
27	10.358	1215	1219	1227	rBV8	10024	15628	0.22%	0.020%
28	10.546	1242	1251	1278	rBV	503470	1807509	25.09%	2.257%
29	10.863	1298	1305	1322	rVV	1577474	3293894	45.73%	4.113%
30	10.987	1324	1326	1330	rVV5	9709	13528	0.19%	0.017%
31	11.057	1331	1338	1365	rVB	270320	983330	13.65%	1.228%
32	11.557	1414	1423	1425	rBV10	10790	25702	0.36%	0.032%
33	11.728	1448	1452	1455	rBV6	7324	13463	0.19%	0.017%
34	11.852	1468	1473	1480	rVB	24448	39927	0.55%	0.050%
35	12.057	1504	1508	1512	rBV6	8571	11998	0.17%	0.015%
36	12.257	1537	1542	1548	rVV3	13910	26568	0.37%	0.033%

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Integrator: RTE
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 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 >
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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	12.481	1574	1580	1584	rVV8	14992	29932	0.42%	0.037%
38	12.563	1587	1594	1601	rVV8	13444	32970	0.46%	0.041%
39	12.757	1620	1627	1635	rBV5	13916	34456	0.48%	0.043%
40	13.034	1669	1674	1676	rBV6	10613	18592	0.26%	0.023%
41	13.193	1695	1701	1708	rBV3	22905	47845	0.66%	0.060%
42	13.428	1738	1741	1747	rVV8	13040	23770	0.33%	0.030%
43	13.487	1747	1751	1755	rVV	25046	41395	0.57%	0.052%
44	13.551	1756	1762	1775	rVV3	139473	295611	4.10%	0.369%
45	13.663	1777	1781	1787	rVB8	9046	15923	0.22%	0.020%
46	13.869	1812	1816	1819	rBV6	7780	15466	0.21%	0.019%
47	13.940	1825	1828	1831	rBV3	25800	33600	0.47%	0.042%
48	13.987	1831	1836	1847	rVV3	282942	433137	6.01%	0.541%
49	14.104	1847	1856	1876	rVV	1478408	2705934	37.57%	3.379%
50	14.387	1897	1904	1918	rBV2	2001588	3445080	47.83%	4.302%
51	14.504	1920	1924	1929	rVB4	34489	49599	0.69%	0.062%
52	14.551	1929	1932	1938	rVB2	41249	61161	0.85%	0.076%
53	14.622	1938	1944	1950	rBV	436993	631321	8.76%	0.788%
54	14.693	1950	1956	1965	rBV	2412825	3764668	52.26%	4.701%
55	14.816	1974	1977	1983	rVB7	34906	46023	0.64%	0.057%
56	14.916	1988	1994	1998	rBV3	42695	72900	1.01%	0.091%
57	15.040	2009	2015	2023	rBV6	72765	252252	3.50%	0.315%
58	15.169	2034	2037	2047	rVB6	26242	64082	0.89%	0.080%
59	15.310	2057	2061	2066	rBV7	15313	29067	0.40%	0.036%
60	15.457	2082	2086	2090	rVB6	17671	23083	0.32%	0.029%
61	15.504	2090	2094	2100	rVV2	33556	55610	0.77%	0.069%
62	15.569	2101	2105	2107	rVV4	13912	23290	0.32%	0.029%
63	15.593	2107	2109	2113	rVB5	17590	19970	0.28%	0.025%
64	15.693	2118	2126	2142	rBV	1969969	3628814	50.38%	4.531%
65	15.887	2153	2159	2172	rBV4	140315	465732	6.47%	0.582%
66	15.998	2175	2178	2184	rVB8	25716	39557	0.55%	0.049%
67	16.063	2184	2189	2197	rBV	241782	489318	6.79%	0.611%
68	16.381	2237	2243	2248	rBV	76473	149380	2.07%	0.187%
69	16.545	2266	2271	2277	rBV3	132285	248650	3.45%	0.310%
70	16.834	2317	2320	2325	rVB6	30230	44071	0.61%	0.055%
71	16.987	2342	2346	2349	rBV6	18138	26118	0.36%	0.033%
72	17.163	2372	2376	2381	rBV3	59558	83393	1.16%	0.104%
73	17.328	2400	2404	2410	rVB3	57215	87837	1.22%	0.110%
74	17.392	2412	2415	2419	rBV5	33555	45421	0.63%	0.057%
75	17.457	2420	2426	2431	rBV	2493964	3749330	52.05%	4.681%
76	17.498	2431	2433	2438	rVV2	298139	443804	6.16%	0.554%

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77	17.563	2438	2444	2464	rVB	1739381	3241618	45.00%	4.048%
78	17.904	2498	2502	2507	rBV2	76444	106414	1.48%	0.133%
79	18.010	2517	2520	2524	rBV3	56724	78865	1.09%	0.098%
80	18.081	2530	2532	2537	rVB5	54981	66110	0.92%	0.083%
81	18.192	2548	2551	2557	rBV4	53036	70367	0.98%	0.088%
82	18.251	2557	2561	2566	rBV	432322	581097	8.07%	0.726%
83	18.310	2566	2571	2579	rVV	2301236	3077964	42.73%	3.843%
84	18.381	2580	2583	2593	rVB	125679	238530	3.31%	0.298%
85	19.181	2716	2719	2721	rBV	71703	83500	1.16%	0.104%
86	19.392	2752	2755	2757	rBV	148705	184116	2.56%	0.230%
87	19.416	2757	2759	2761	rVV	307165	361211	5.01%	0.451%
88	19.457	2764	2766	2774	rVB	630048	847321	11.76%	1.058%
89	19.533	2775	2779	2790	rBV	803914	1152029	15.99%	1.438%
90	19.786	2817	2822	2826	rBV	2715078	4074394	56.56%	5.087%
91	20.251	2899	2901	2907	rBV	138985	223026	3.10%	0.278%
92	21.192	3058	3061	3064	rBV4	347413	412339	5.72%	0.515%
93	21.228	3064	3067	3073	rVV2	726262	1102972	15.31%	1.377%
94	21.398	3093	3096	3100	rBV	620893	697837	9.69%	0.871%
95	21.545	3116	3121	3126	rBV2	2572632	3902907	54.18%	4.873%
96	22.110	3214	3217	3221	rVB	665787	817973	11.36%	1.021%
97	23.210	3399	3404	3413	rVB	2657064	4478843	62.18%	5.592%
98	23.921	3520	3525	3531	rVB2	1516551	2775731	38.54%	3.466%
99	24.092	3548	3554	3562	rVB	1617917	3522901	48.91%	4.399%
100	24.639	3641	3647	3656	rVB	3347671	7203139	100.00%	8.994%

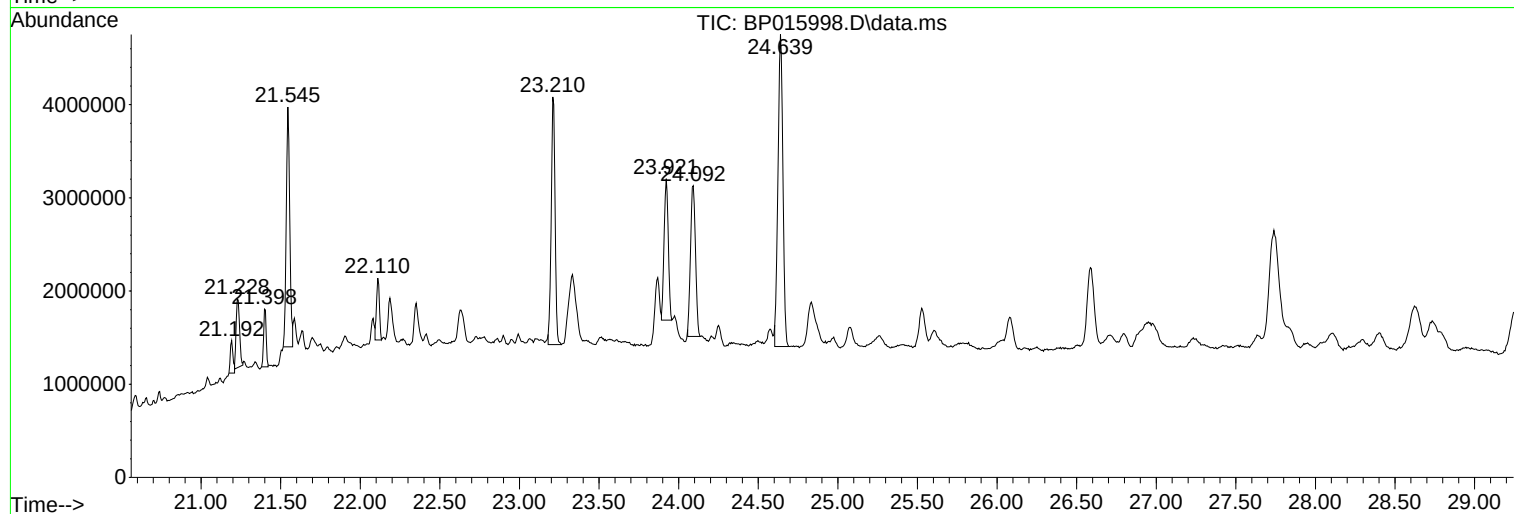
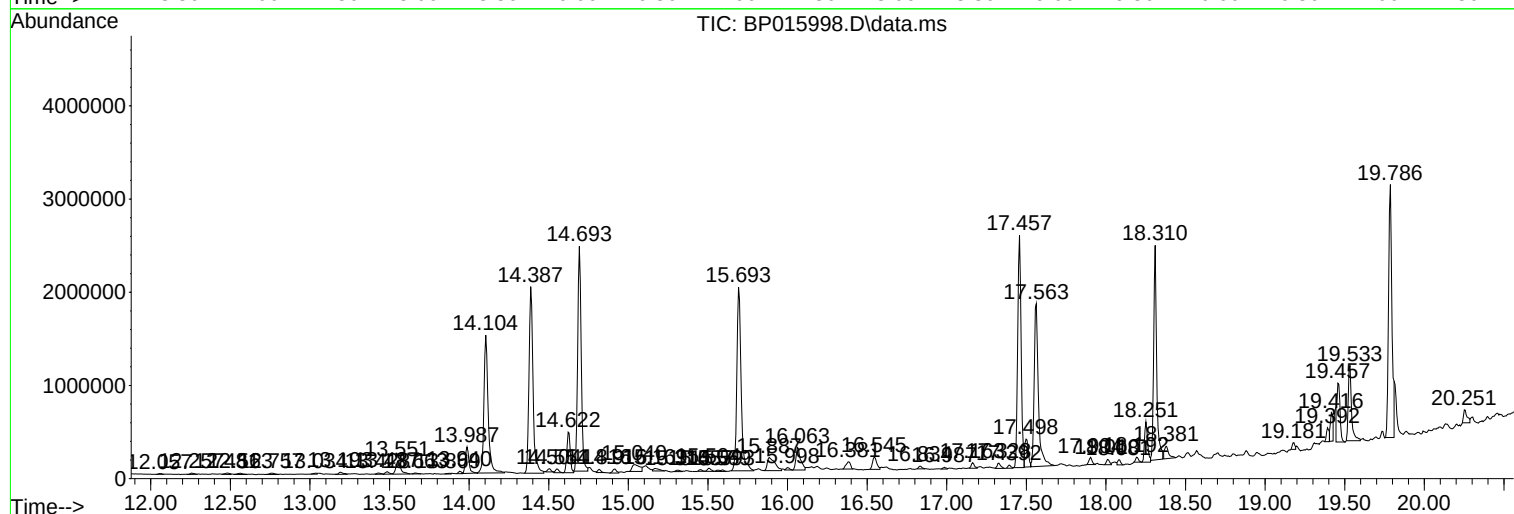
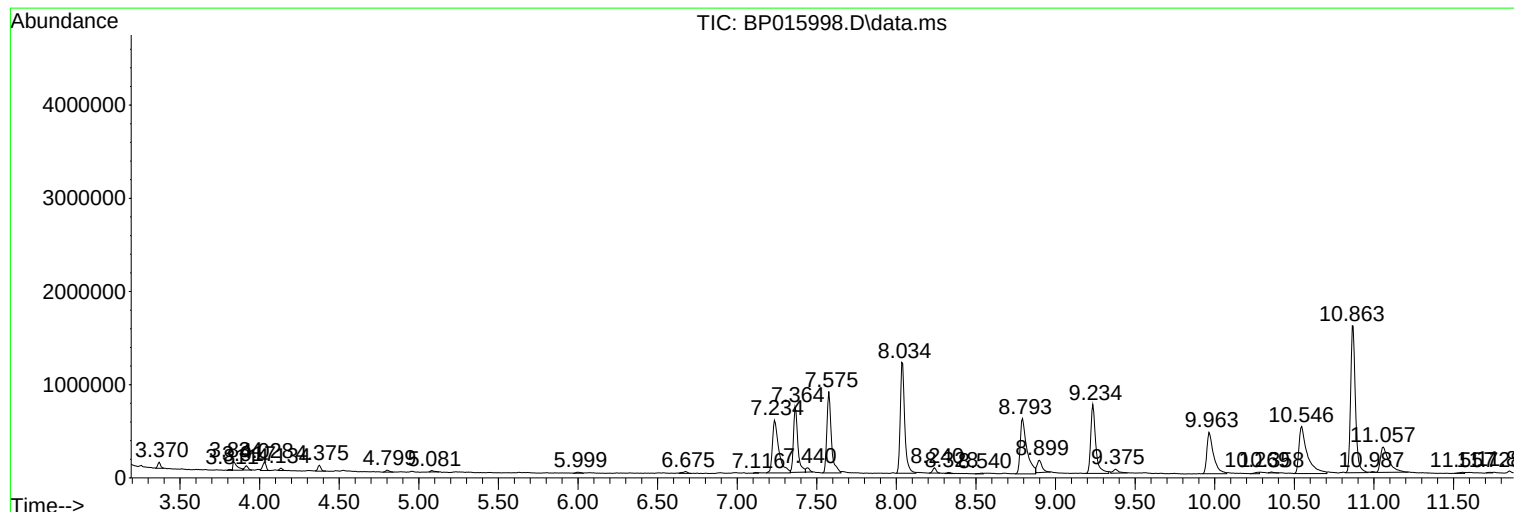
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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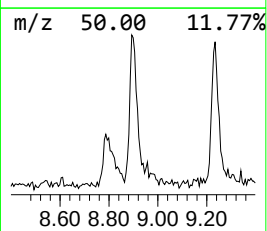
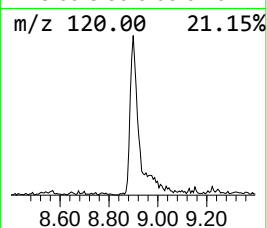
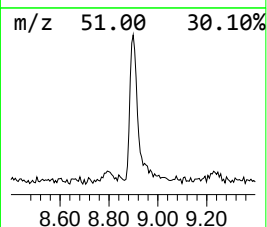
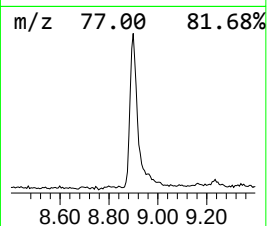
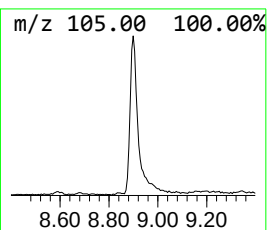
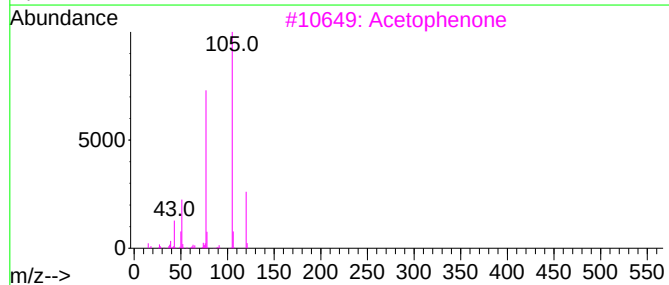
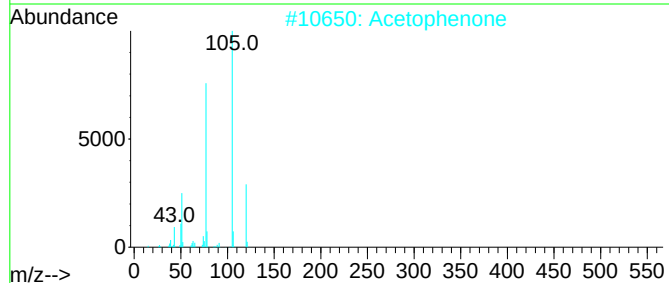
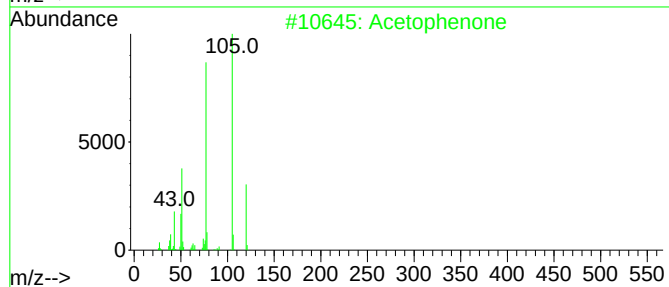
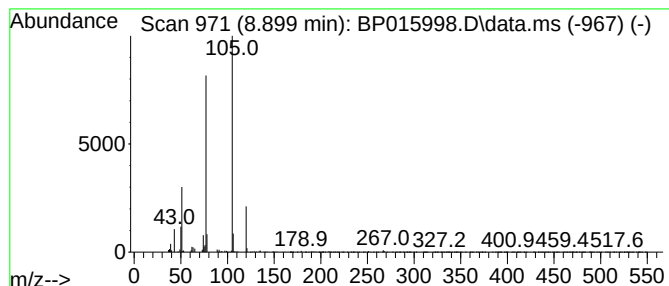
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.899	2.49 ng/ul	285539	1,4-Dichlorobenzene-d4	8.040

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	95
2			Acetophenone	120	C8H8O	000098-86-2	94
3			Acetophenone	120	C8H8O	000098-86-2	91
4			Acetophenone	120	C8H8O	000098-86-2	91
5			Acetophenone	120	C8H8O	000098-86-2	87



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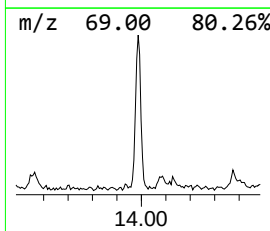
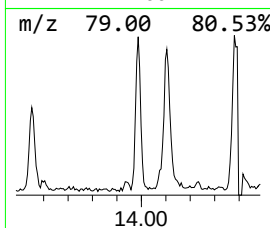
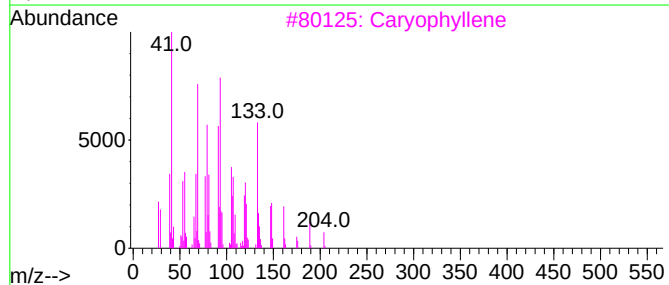
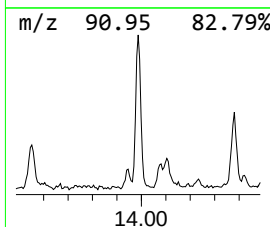
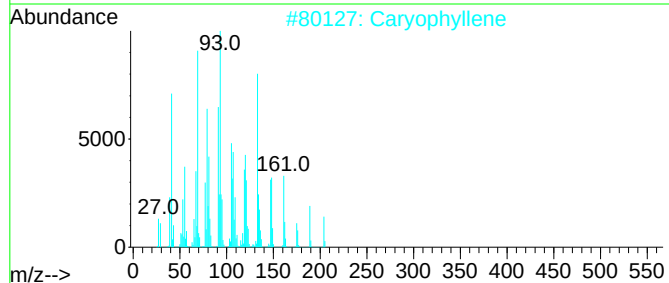
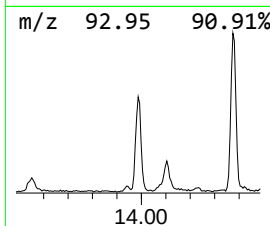
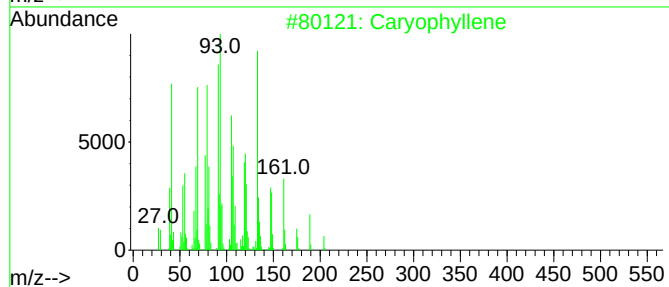
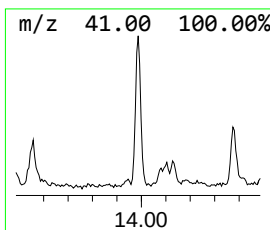
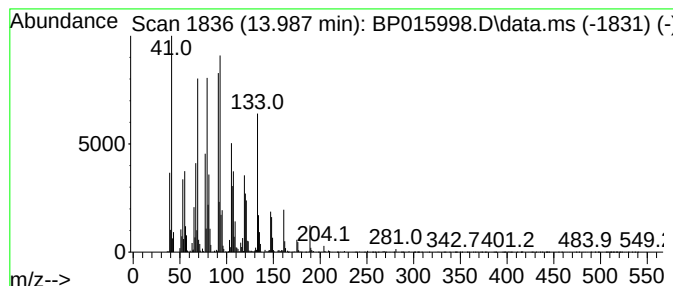
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Caryophyllene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.987	2.30 ng/ul	433137	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Caryophyllene	204	C15H24	000087-44-5	99
2			Caryophyllene	204	C15H24	000087-44-5	98
3			Caryophyllene	204	C15H24	000087-44-5	98
4			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	000118-65-0	95
5			Bicyclo[7.2.0]undec-4-ene, 4,11,...	204	C15H24	013877-93-5	94



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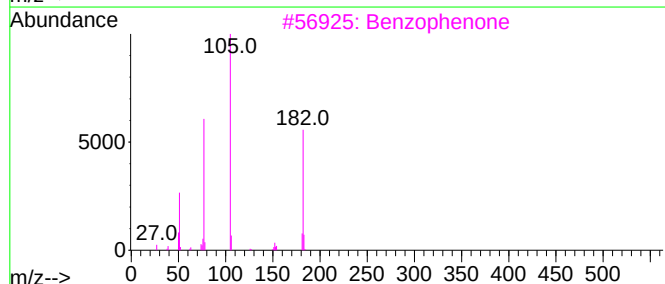
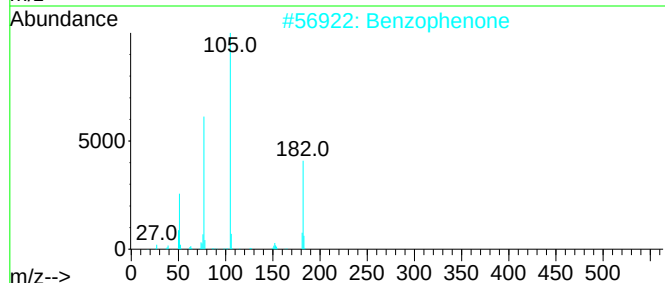
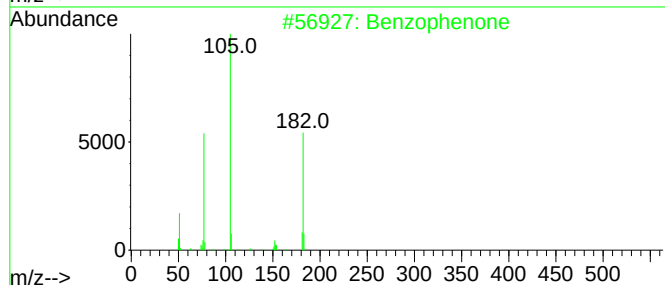
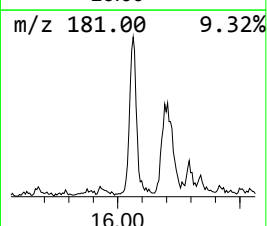
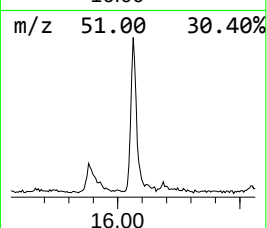
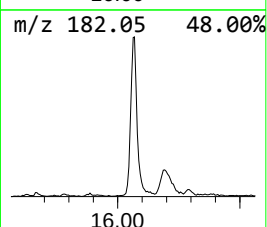
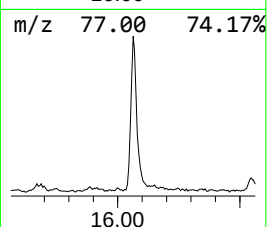
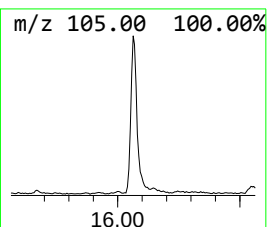
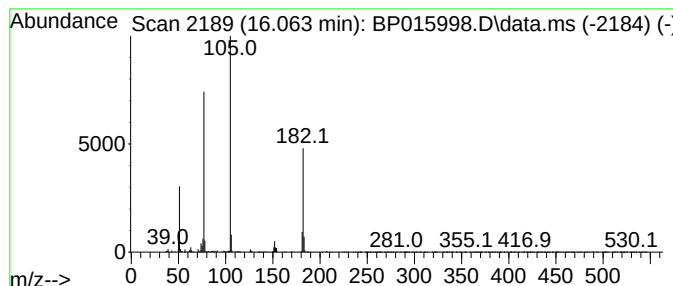
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzophenone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.063	2.60 ng/ul	489318	Acenaphthene-d10	14.693

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	76



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Data File : BP015998.D
Acq On : 05 Jul 2023 04:07
Operator : MA/JU
Sample : 03244-22
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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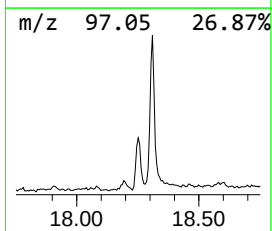
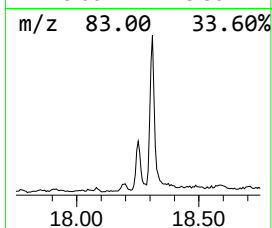
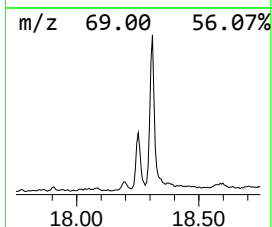
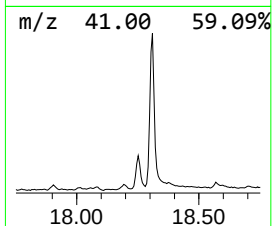
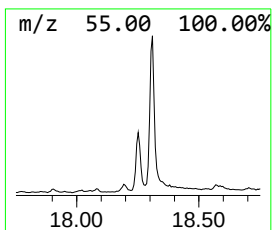
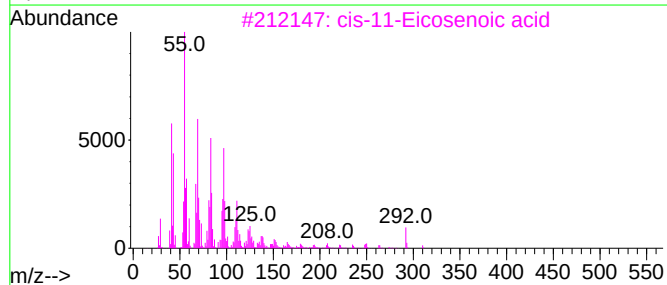
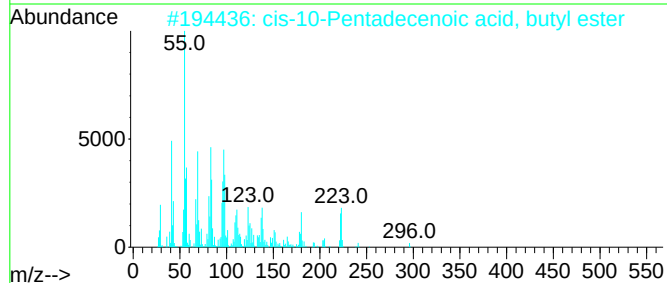
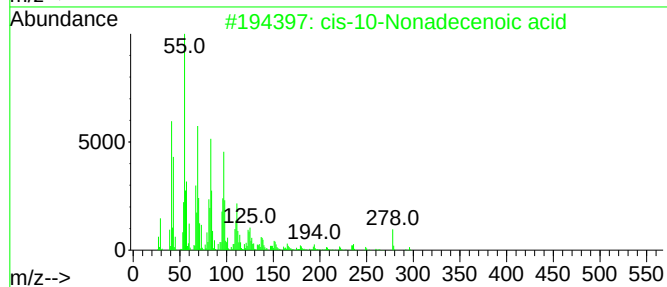
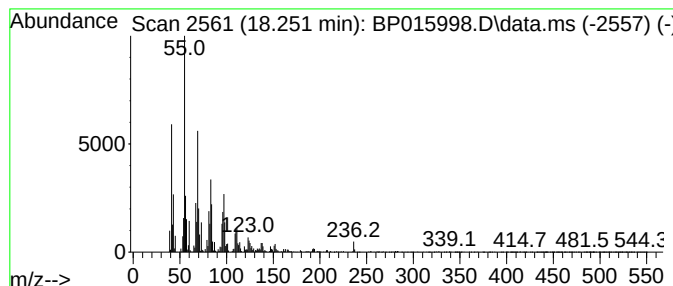
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 cis-10-Nonadecenoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.251	3.10 ng/ul	581097	Phenanthrene-d10	17.457

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-10-Nonadecenoic acid	296	C19H36O2	073033-09-7	98
2			cis-10-Pentadecenoic acid, butyl...	296	C19H36O2	1000405-17-6	95
3			cis-11-Eicosenoic acid	310	C20H38O2	005561-99-9	91
4			cis-Vaccenic acid	282	C18H34O2	000506-17-2	91
5			Palmitoleic acid	254	C16H30O2	000373-49-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
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Sample : 03244-22
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ALS Vial : 44 Sample Multiplier: 1

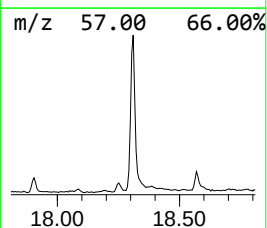
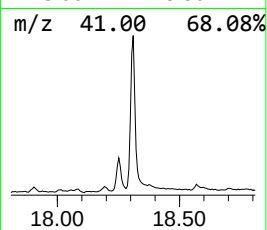
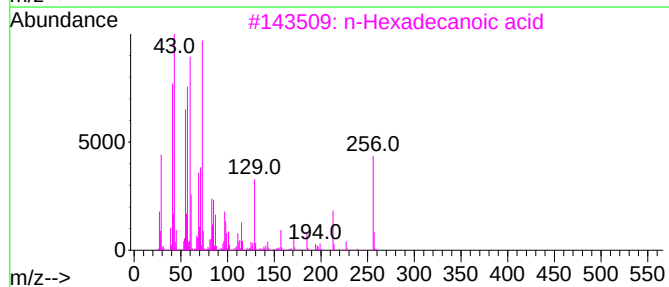
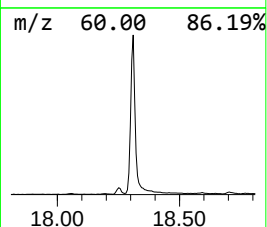
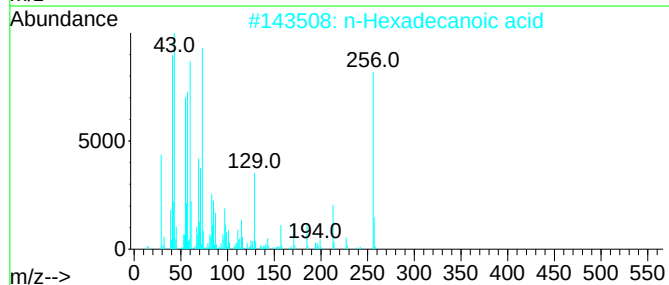
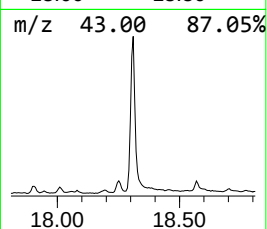
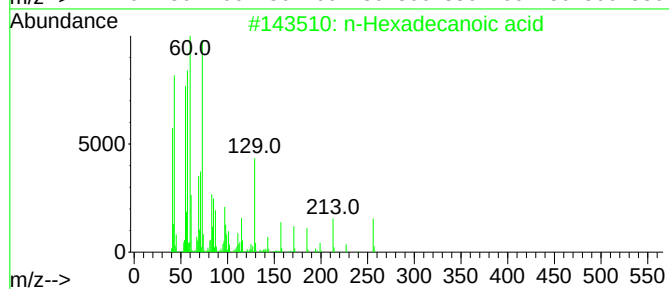
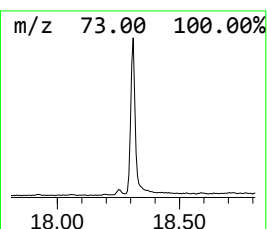
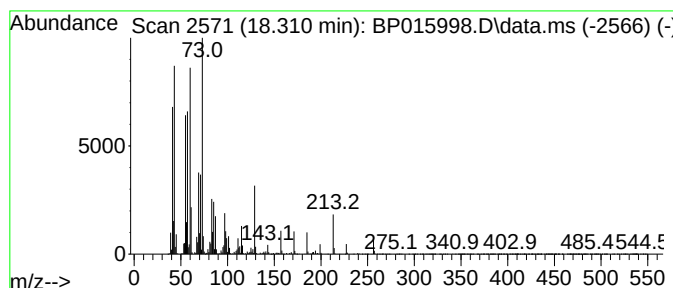
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.310	16.42 ng/ul	3077960	Phenanthrene-d10	17.457	
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	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015998.D
Acq On : 05 Jul 2023 04:07
Operator : MA/JU
Sample : 03244-22
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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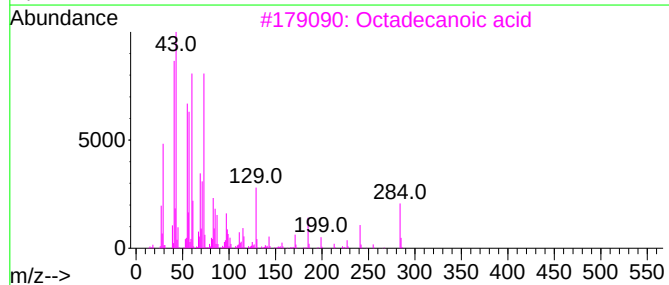
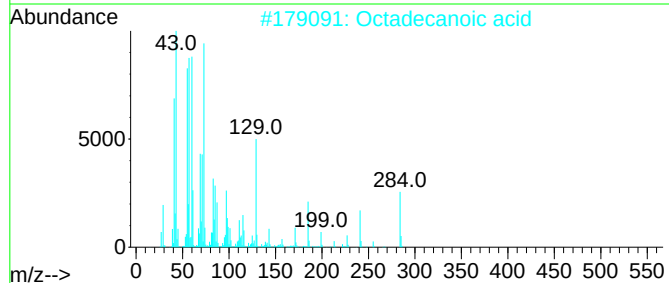
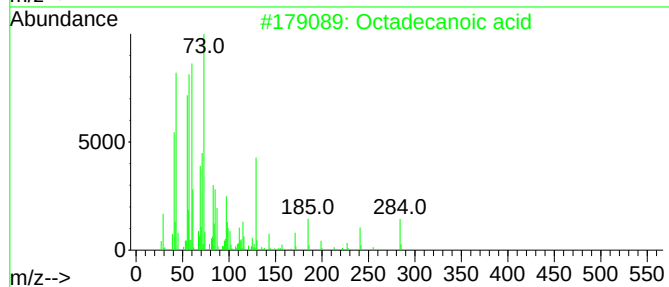
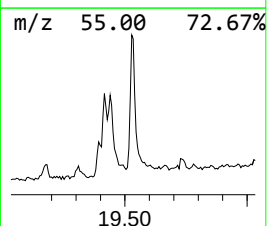
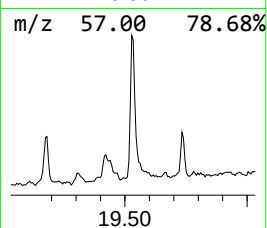
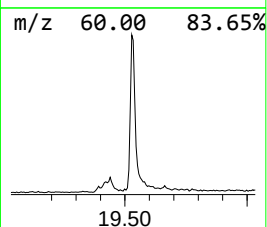
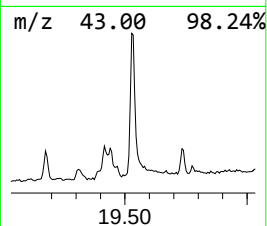
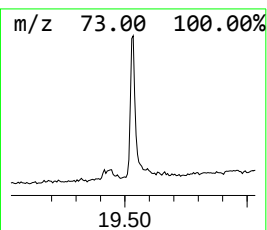
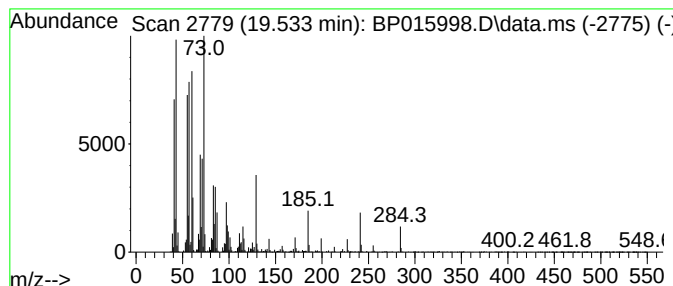
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.534	5.90 ng/ul	1152030	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015998.D
Acq On : 05 Jul 2023 04:07
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Sample : 03244-22
Misc :
ALS Vial : 44 Sample Multiplier: 1

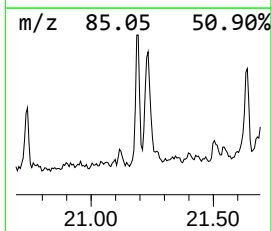
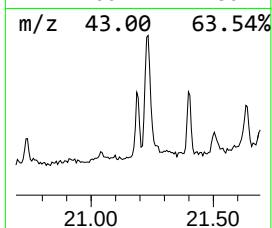
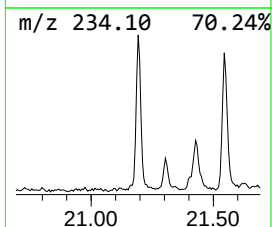
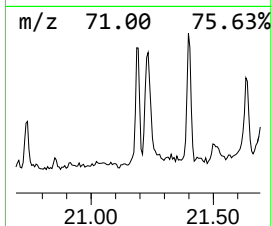
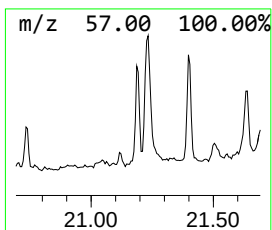
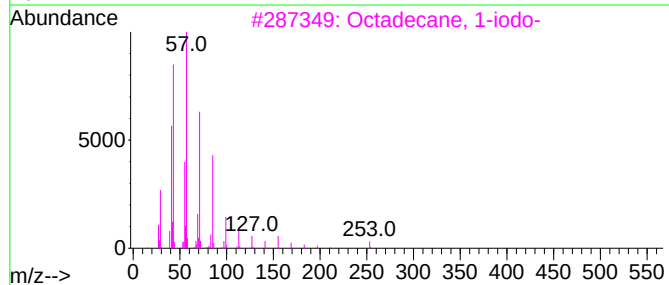
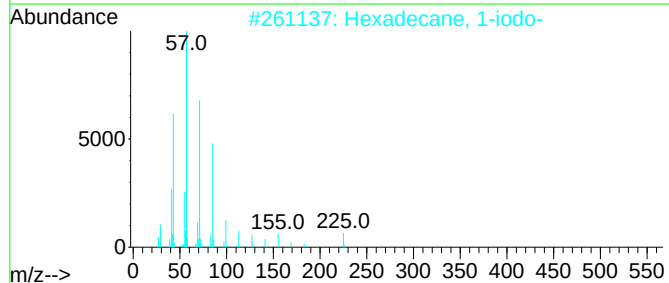
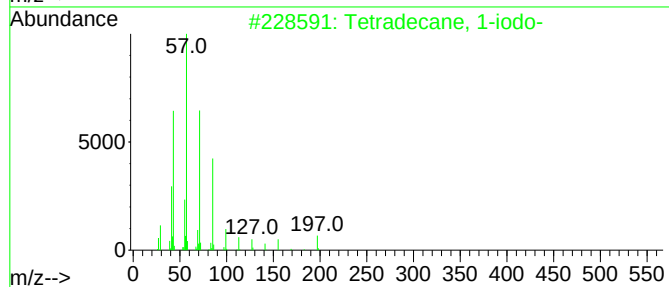
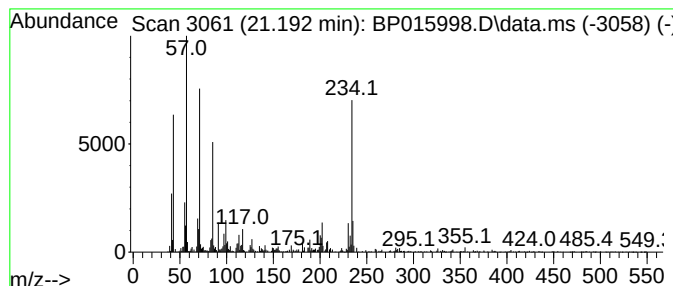
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Tetradecane, 1-iodo- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.192	2.11 ng/ul	412339	Chrysene-d12	21.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	53
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	53
3	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	49
4	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	49
5	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
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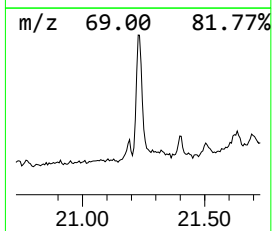
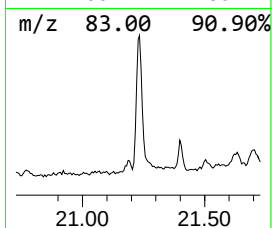
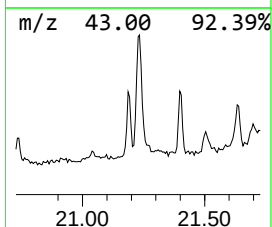
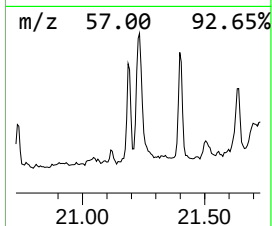
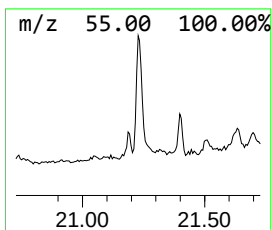
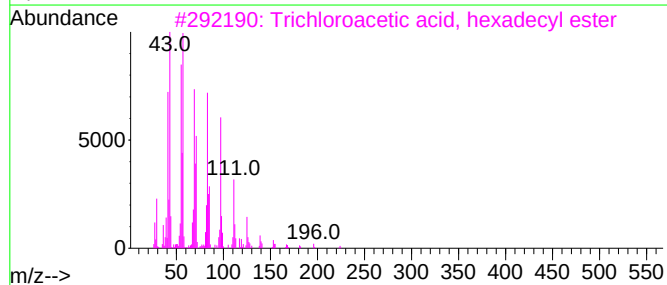
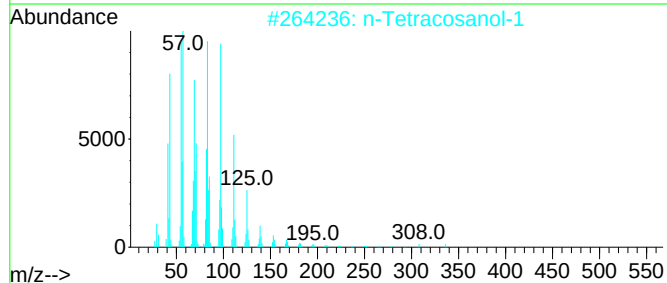
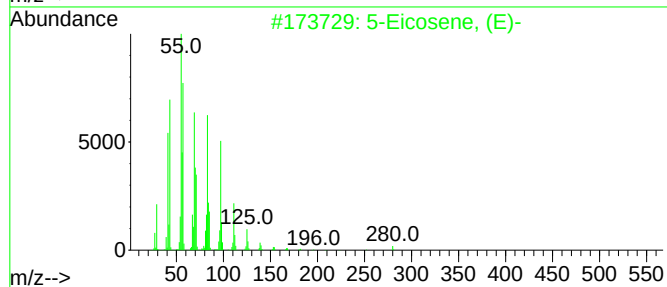
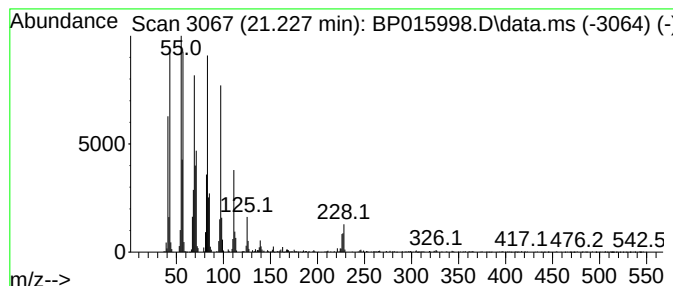
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.227	5.65 ng/ul	1102970	Chrysene-d12	21.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	99
2	n-Tetracosanol-1	354	C24H50O	000506-51-4	94
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
4	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5	Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
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Acq On : 05 Jul 2023 04:07
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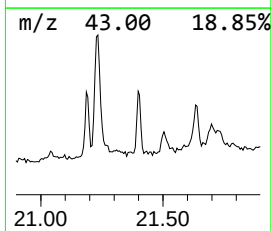
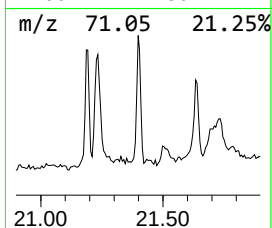
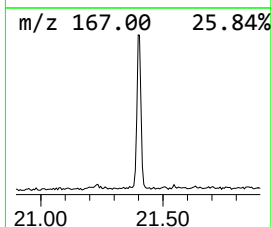
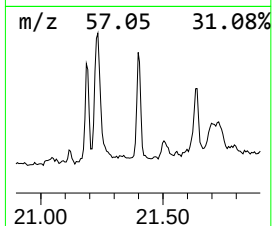
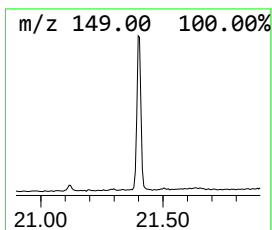
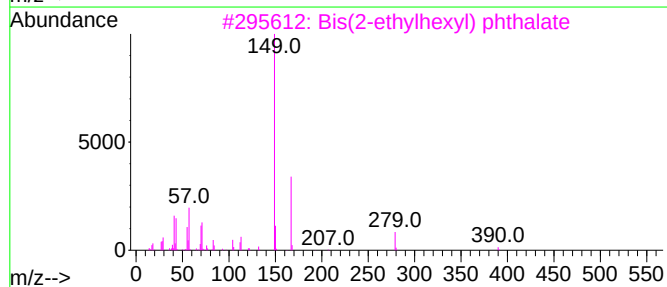
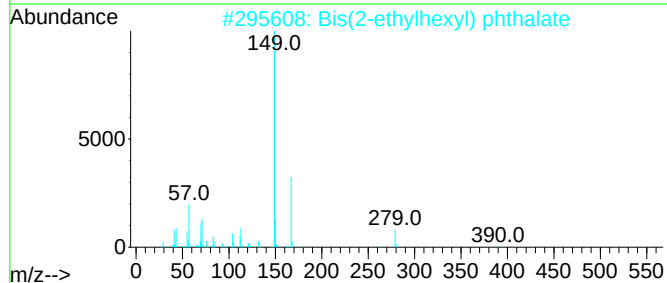
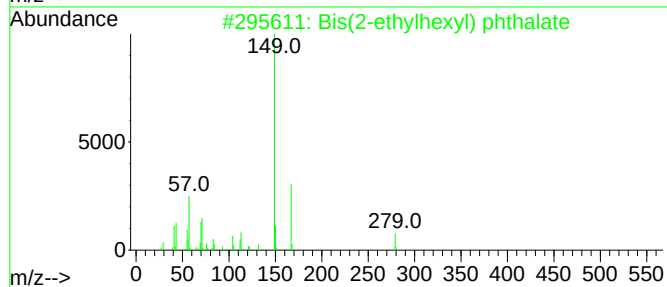
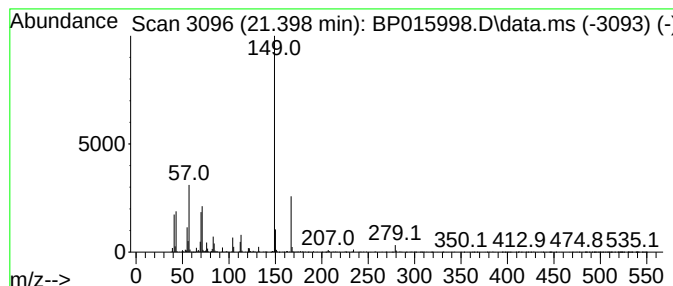
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Bis(2-ethylhexyl) phthalate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.398	3.58 ng/ul	697837	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	87
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	86
4			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	72
5			Phthalic acid, isohexyl 2-methyl...	334	C20H30O4	1000356-91-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015998.D
Acq On : 05 Jul 2023 04:07
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Sample : 03244-22
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ALS Vial : 44 Sample Multiplier: 1

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ClientSampleId :
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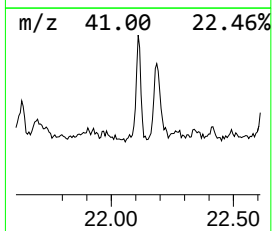
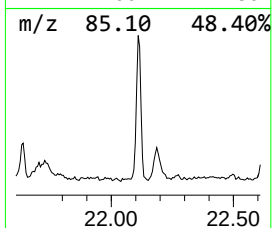
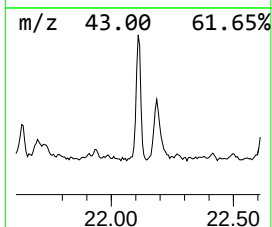
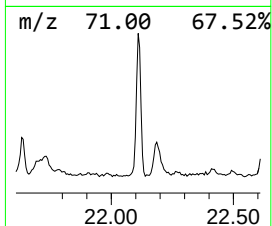
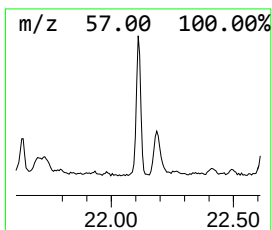
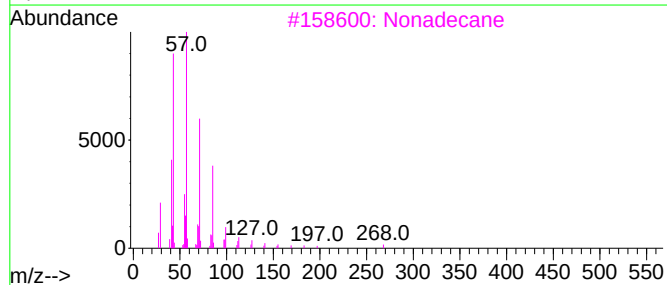
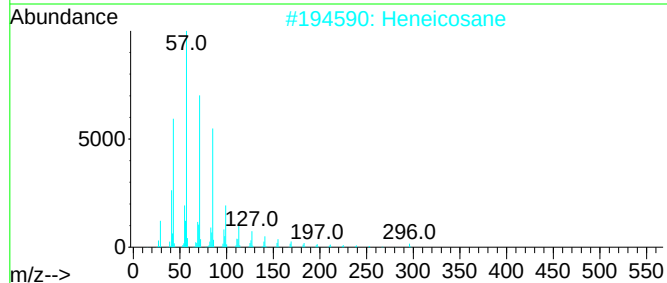
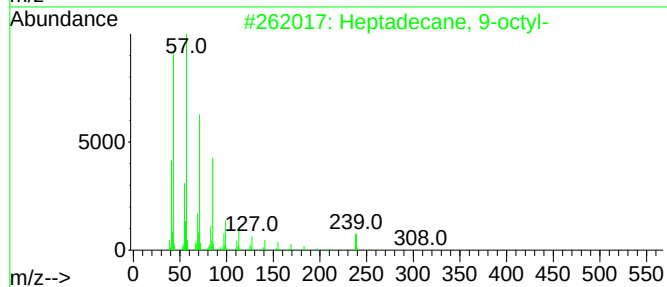
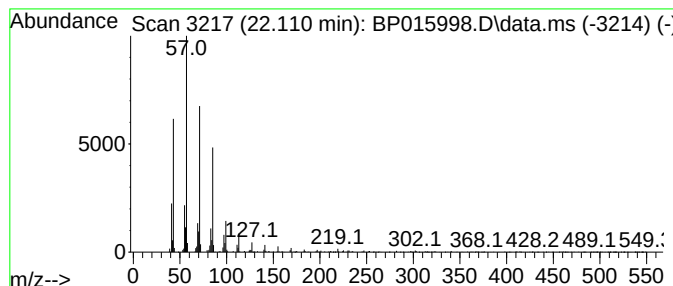
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.110	4.19 ng/ul	817973	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
2			Heneicosane	296	C21H44	000629-94-7	91
3			Nonadecane	268	C19H40	000629-92-5	91
4			Octadecane	254	C18H38	000593-45-3	91
5			Tricosane	324	C23H48	000638-67-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015998.D
Acq On : 05 Jul 2023 04:07
Operator : MA/JU
Sample : 03244-22
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGG1

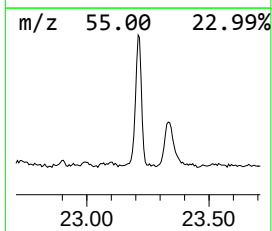
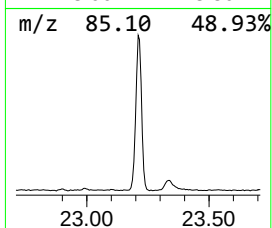
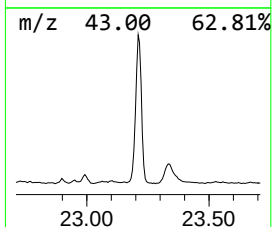
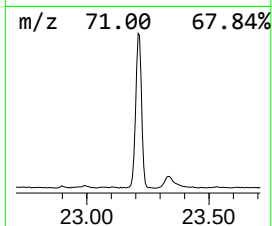
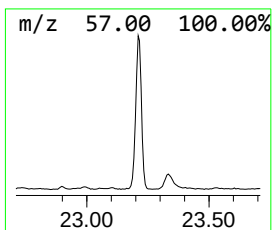
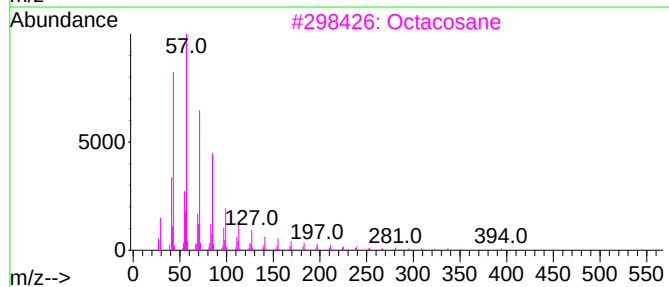
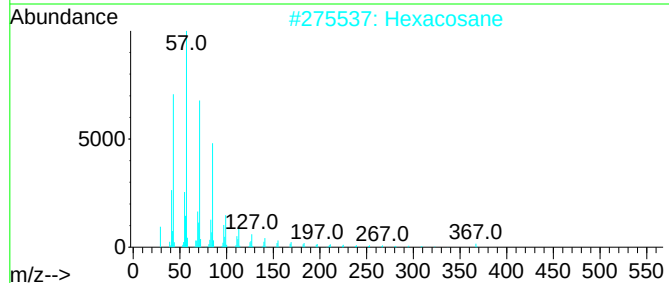
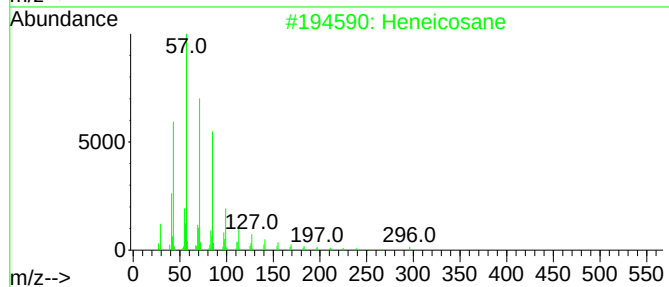
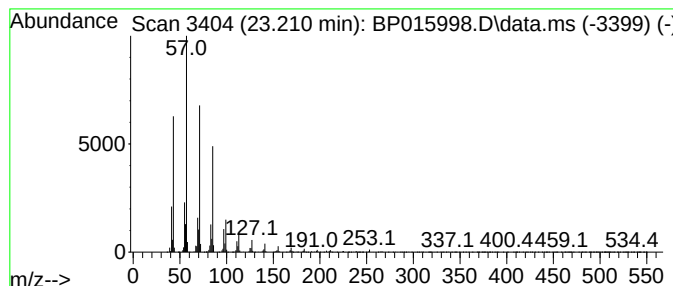
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.210	25.43 ng/ul	4478840	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	97
2		Hexacosane	366	C26H54	000630-01-3	91
3		Octacosane	394	C28H58	000630-02-4	91
4		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
Data File : BP015998.D
Acq On : 05 Jul 2023 04:07
Operator : MA/JU
Sample : 03244-22
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGG1

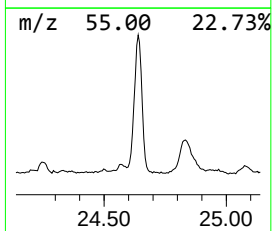
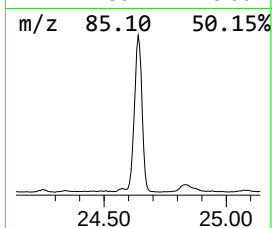
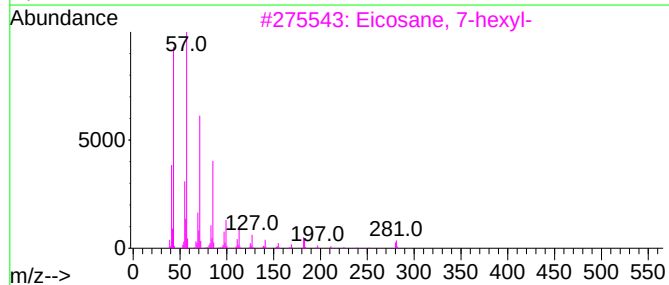
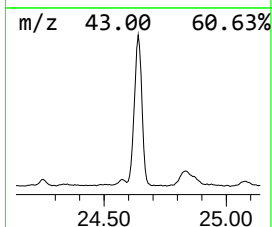
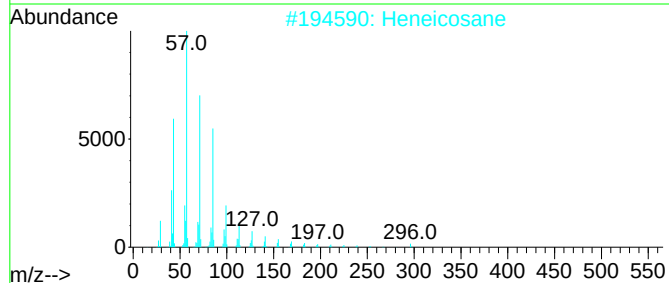
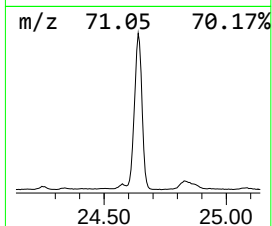
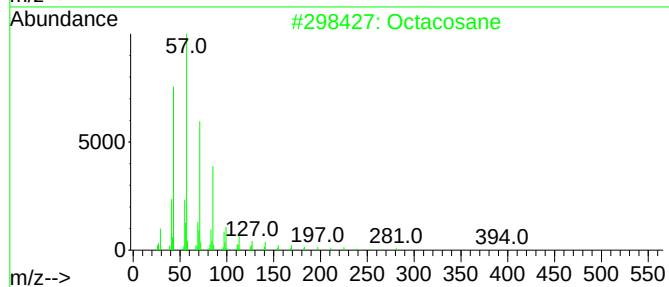
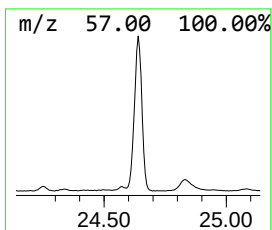
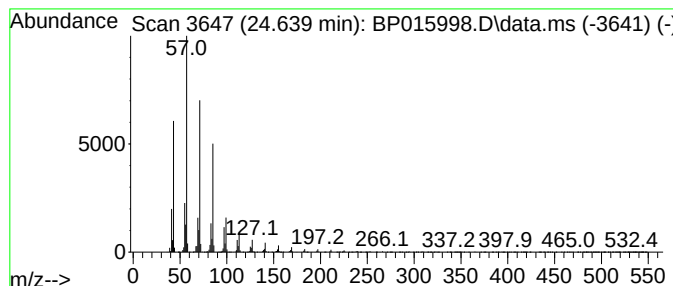
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.639	40.89 ng/ul	7203140	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	91
2		Heneicosane	296	C21H44	000629-94-7	91
3		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4		Triacontane	422	C30H62	000638-68-6	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070423\
 Data File : BP015998.D
 Acq On : 05 Jul 2023 04:07
 Operator : MA/JU
 Sample : 03244-22
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGG1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Aceto phenone	8.899	2.5	ng/ul	285539	1	8.040	2293720	20.0
Caryophyllene	13.987	2.3	ng/ul	433137	3	14.693	3764670	20.0
(1R,2S,6S,7S,8S...	14.622	3.4	ng/ul	631321	3	14.693	3764670	20.0
Benzophenone	16.063	2.6	ng/ul	489318	3	14.693	3764670	20.0
cis-10-Nonadece...	18.251	3.1	ng/ul	581097	4	17.457	3749330	20.0
n-Hexadecanoic ...	18.310	16.4	ng/ul	3077960	4	17.457	3749330	20.0
Octadecanoic acid	19.534	5.9	ng/ul	1152030	5	21.545	3902910	20.0
Tetradecane, 1-...	21.192	2.1	ng/ul	412339	5	21.545	3902910	20.0
(DEL) Alkane: S...	21.227	5.7	ng/ul	1102970	5	21.545	3902910	20.0
Bis(2-ethylhexy...	21.398	3.6	ng/ul	697837	5	21.545	3902910	20.0
(DEL) Alkane: S...	22.110	4.2	ng/ul	817973	5	21.545	3902910	20.0
(DEL) Alkane: S...	23.210	25.4	ng/ul	4478840	6	24.092	3522900	20.0
(DEL) Alkane: S...	24.639	40.9	ng/ul	7203140	6	24.092	3522900	20.0