

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
 Data File : BP015670.D
 Acq On : 23 Jun 2023 14:23
 Operator : MA/JU
 Sample : 03084-19
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFN6

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-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015670.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	33	36	44	rVB	18478	26418	1.30%	0.114%
2	3.846	110	112	117	rVB4	5403	5417	0.27%	0.023%
3	3.922	122	125	126	rBV2	4527	5481	0.27%	0.024%
4	3.952	126	130	136	rVB	29389	39474	1.95%	0.170%
5	4.069	146	150	155	rVB	16226	24573	1.21%	0.106%
6	4.175	164	168	173	rVB2	10456	16970	0.84%	0.073%
7	4.564	231	234	237	rVB5	4127	4735	0.23%	0.020%
8	4.752	263	266	270	rBV6	6576	7516	0.37%	0.032%
9	5.122	327	329	333	rVB5	4073	4643	0.23%	0.020%
10	5.887	456	459	463	rBV6	2632	4690	0.23%	0.020%
11	6.040	479	485	492	rBV6	5396	12200	0.60%	0.053%
12	6.311	526	531	536	rBV8	4752	10934	0.54%	0.047%
13	6.928	631	636	637	rBV5	3299	5272	0.26%	0.023%
14	7.246	683	690	711	rBV	133052	316541	15.64%	1.366%
15	7.405	711	717	730	rVB	144707	254552	12.57%	1.098%
16	7.610	746	752	760	rBV	198437	336573	16.62%	1.452%
17	7.893	798	800	806	rVB7	2855	4838	0.24%	0.021%
18	8.081	825	832	846	rBV	503861	852400	42.10%	3.678%
19	8.816	949	957	970	rBV	80073	193692	9.57%	0.836%
20	8.928	970	976	986	rVB	44578	94627	4.67%	0.408%
21	9.263	1025	1033	1050	rBV	161815	346581	17.12%	1.495%
22	9.628	1091	1095	1099	rVB6	3319	4847	0.24%	0.021%
23	9.993	1148	1157	1166	rBV	138933	282037	13.93%	1.217%
24	10.257	1194	1202	1203	rVV8	4540	8998	0.44%	0.039%
25	10.304	1207	1210	1215	rVV7	3648	6021	0.30%	0.026%
26	10.375	1215	1222	1228	rVV7	3747	10074	0.50%	0.043%
27	10.557	1245	1253	1272	rBV	116749	344975	17.04%	1.488%
28	10.910	1305	1313	1331	rVV	600360	1098574	54.26%	4.740%
29	11.098	1338	1345	1354	rBV2	16855	48668	2.40%	0.210%
30	11.363	1384	1390	1391	rBV6	3722	6424	0.32%	0.028%
31	11.604	1427	1431	1436	rVB7	2965	4332	0.21%	0.019%
32	11.769	1453	1459	1462	rBV8	3492	7570	0.37%	0.033%
33	11.904	1478	1482	1487	rBV5	6284	9713	0.48%	0.042%
34	12.316	1546	1552	1555	rBV6	4055	8030	0.40%	0.035%
35	12.469	1572	1578	1588	rVV4	59292	110365	5.45%	0.476%
36	12.587	1592	1598	1603	rBV8	5065	10187	0.50%	0.044%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	12.787	1626	1632	1639	rBV9	4301	9439	0.47%	0.041%
38	13.245	1706	1710	1717	rVB8	5461	10453	0.52%	0.045%
39	13.534	1755	1759	1762	rBV5	5359	8422	0.42%	0.036%
40	13.610	1768	1772	1778	rVB3	14619	24049	1.19%	0.104%

41	14.051	1842	1847	1852	rBV7	7260	14306	0.71%	0.062%
42	14.140	1852	1862	1875	rBV	339356	582312	28.76%	2.513%
43	14.287	1884	1887	1891	rBV6	3191	4643	0.23%	0.020%
44	14.428	1903	1911	1926	rBV	424834	715416	35.34%	3.087%
45	14.545	1928	1931	1935	rVB5	4609	7298	0.36%	0.031%

46	14.604	1938	1941	1947	rVB7	8492	11658	0.58%	0.050%
47	14.681	1949	1954	1955	rBV5	4384	4968	0.25%	0.021%
48	14.728	1956	1962	1971	rBV	794829	1222184	60.37%	5.273%
49	15.004	2003	2009	2017	rBV4	25862	84730	4.19%	0.366%
50	15.139	2028	2032	2038	rBV8	8145	13882	0.69%	0.060%

51	15.351	2065	2068	2074	rBV8	4241	7660	0.38%	0.033%
52	15.434	2080	2082	2087	rVB6	3320	5162	0.25%	0.022%
53	15.510	2087	2095	2098	rBV9	8034	15640	0.77%	0.067%
54	15.557	2098	2103	2108	rVB6	9199	13837	0.68%	0.060%
55	15.728	2125	2132	2146	rBV	475266	781455	38.60%	3.372%

56	15.881	2152	2158	2168	rBV2	72327	153094	7.56%	0.661%
57	15.951	2168	2170	2176	rVB7	6409	10847	0.54%	0.047%
58	16.092	2188	2194	2203	rBV	55202	98656	4.87%	0.426%
59	16.269	2222	2224	2231	rVB8	3251	5694	0.28%	0.025%
60	16.439	2244	2253	2260	rBV4	20282	54665	2.70%	0.236%

61	16.581	2272	2277	2282	rVV9	8679	16992	0.84%	0.073%
62	16.639	2284	2287	2291	rVB6	3403	6674	0.33%	0.029%
63	16.822	2317	2318	2322	rVB4	5549	4477	0.22%	0.019%
64	16.869	2323	2326	2332	rVB5	13548	19854	0.98%	0.086%
65	17.022	2350	2352	2356	rVB5	5061	4802	0.24%	0.021%

66	17.163	2371	2376	2381	rVB8	7011	14526	0.72%	0.063%
67	17.216	2381	2385	2390	rBV3	16426	26749	1.32%	0.115%
68	17.310	2399	2401	2406	rVB3	5683	5947	0.29%	0.026%
69	17.369	2407	2411	2418	rBV8	17133	26737	1.32%	0.115%
70	17.433	2418	2422	2426	rBV5	14696	22002	1.09%	0.095%

71	17.492	2426	2432	2437	rBV	916028	1342679	66.32%	5.793%
72	17.533	2437	2439	2444	rVV	108724	144176	7.12%	0.622%
73	17.592	2444	2449	2463	rVB	463029	781252	38.59%	3.371%
74	17.910	2500	2503	2508	rVB3	12846	22229	1.10%	0.096%
75	17.951	2508	2510	2513	rBV3	11901	12796	0.63%	0.055%

76	18.098	2531	2535	2538	rBV6	7388	11540	0.57%	0.050%
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77	18.233	2553	2558	2564	rBV9	22798	41610	2.06%	0.180%
78	18.292	2565	2568	2572	rBV5	20987	26702	1.32%	0.115%
79	18.351	2573	2578	2584	rBV2	268491	417597	20.63%	1.802%
80	18.422	2588	2590	2597	rVB	58630	69326	3.42%	0.299%
81	18.539	2606	2610	2615	rBV5	26650	58586	2.89%	0.253%
82	18.622	2621	2624	2630	rBV5	15845	28227	1.39%	0.122%
83	18.992	2684	2687	2693	rVB5	28223	40489	2.00%	0.175%
84	19.233	2724	2728	2731	rBV3	47197	61347	3.03%	0.265%
85	19.392	2752	2755	2760	rVB4	40195	57155	2.82%	0.247%
86	19.492	2768	2772	2779	rVB	454163	625704	30.91%	2.700%
87	19.580	2783	2787	2793	rBV6	69649	106742	5.27%	0.461%
88	19.822	2823	2828	2831	rBV	743681	1064249	52.57%	4.592%
89	19.851	2831	2833	2839	rVB	368514	397743	19.65%	1.716%
90	20.304	2908	2910	2913	rBV4	47886	51327	2.54%	0.221%
91	21.251	3068	3071	3077	rVB2	318506	488279	24.12%	2.107%
92	21.457	3102	3106	3112	rBV	444933	528073	26.08%	2.278%
93	21.586	3121	3128	3132	rBV2	1223171	2024554	100.00%	8.735%
94	21.621	3132	3134	3139	rVB	237522	300482	14.84%	1.297%
95	22.857	3341	3344	3349	rVB	237760	323760	15.99%	1.397%
96	23.298	3415	3419	3424	rVB	404204	635306	31.38%	2.741%
97	23.368	3425	3431	3443	rVB2	366864	1144949	56.55%	4.940%
98	23.974	3529	3534	3540	rBV2	532525	1002283	49.51%	4.325%
99	24.145	3558	3563	3570	rVB	818845	1597593	78.91%	6.893%
100	24.751	3661	3666	3677	rVB	548220	1231439	60.83%	5.313%

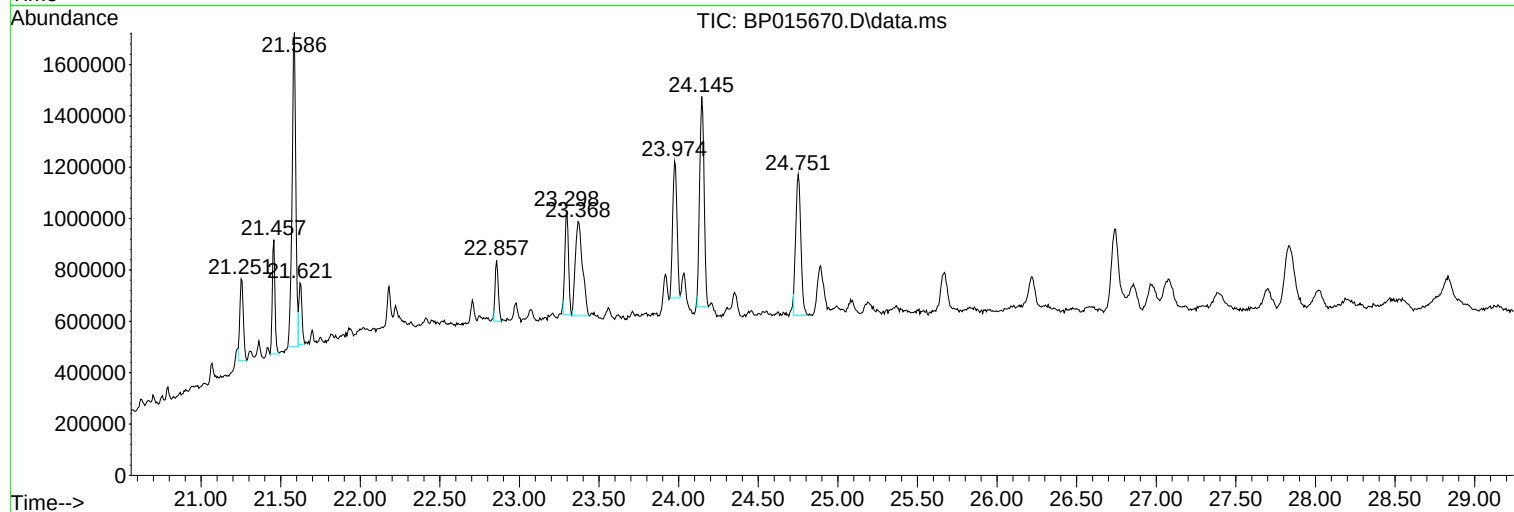
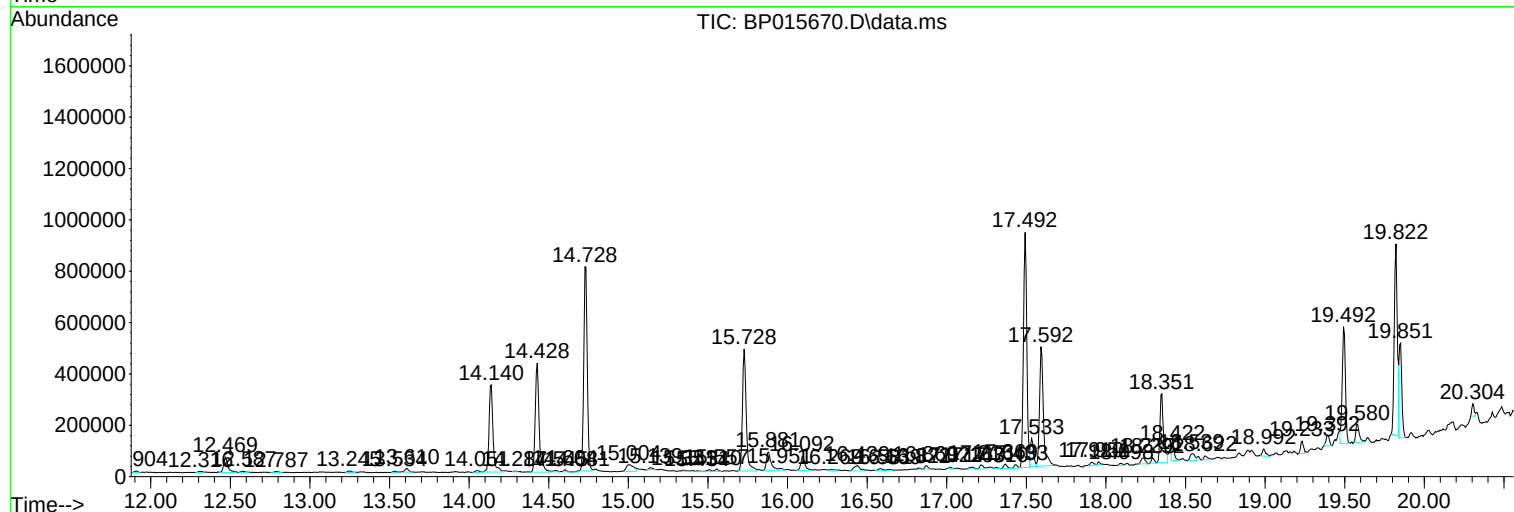
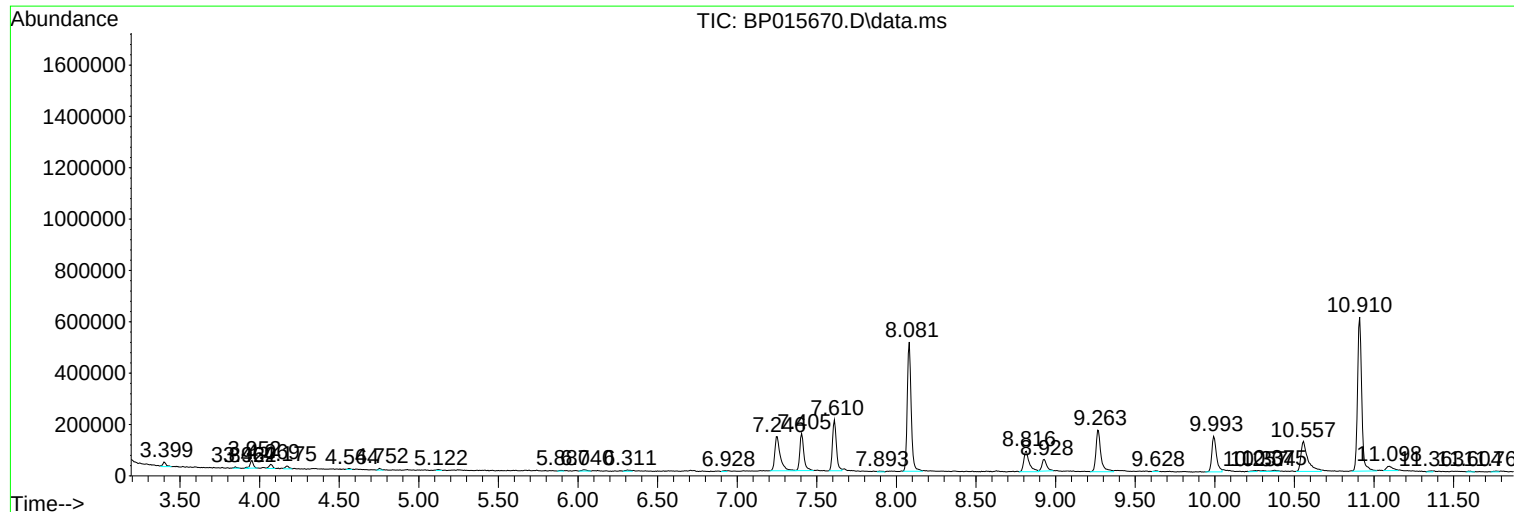
Sum of corrected areas: 23176366

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Misc :
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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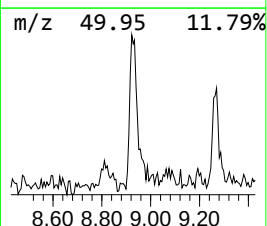
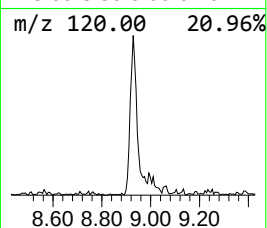
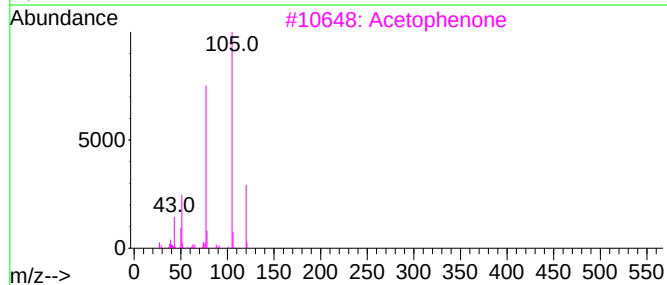
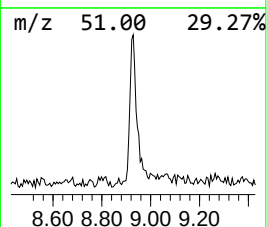
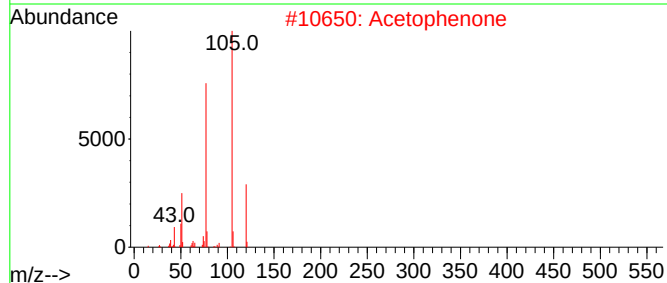
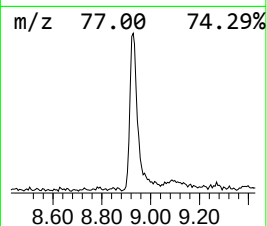
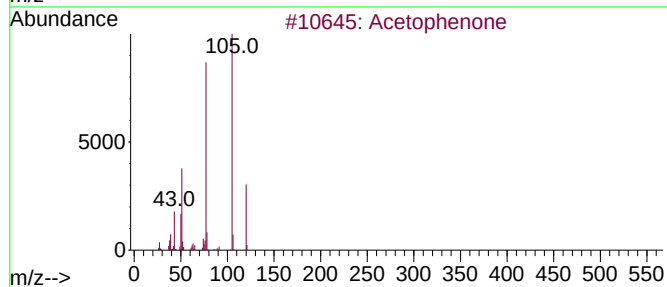
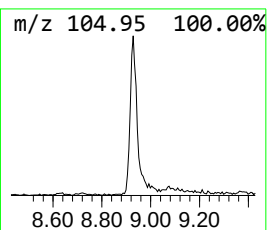
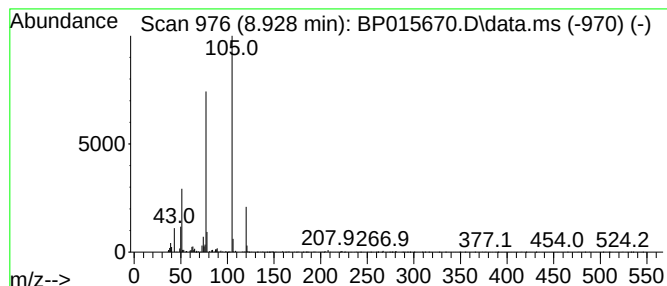
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	2.22 ng/ul	94627	1,4-Dichlorobenzene-d4	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	94
2			Acetophenone	120	C8H8O	000098-86-2	91
3			Acetophenone	120	C8H8O	000098-86-2	91
4			Acetophenone	120	C8H8O	000098-86-2	91
5			Acetophenone	120	C8H8O	000098-86-2	90



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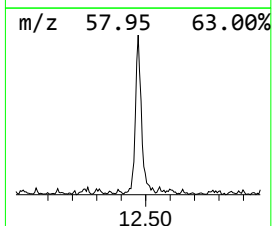
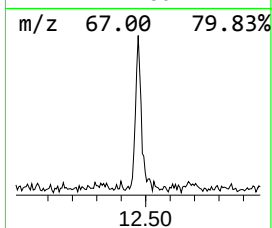
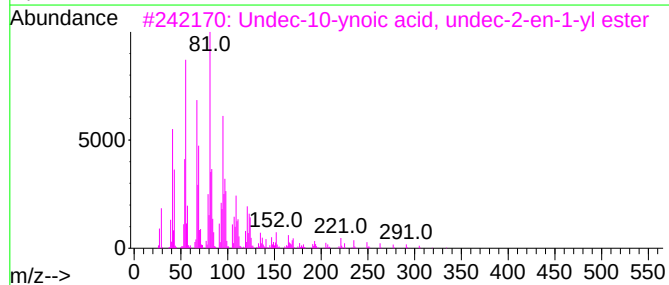
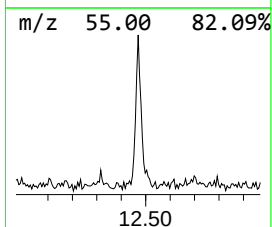
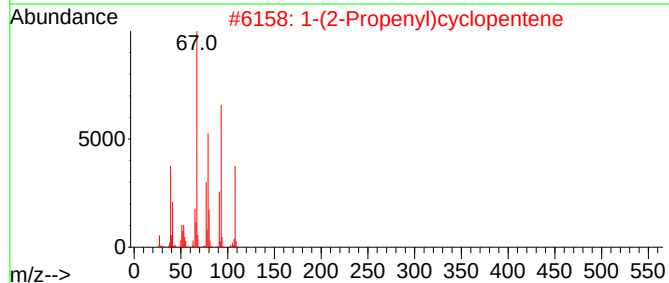
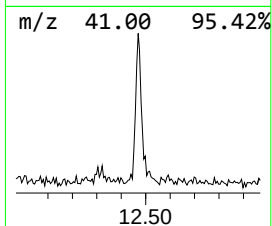
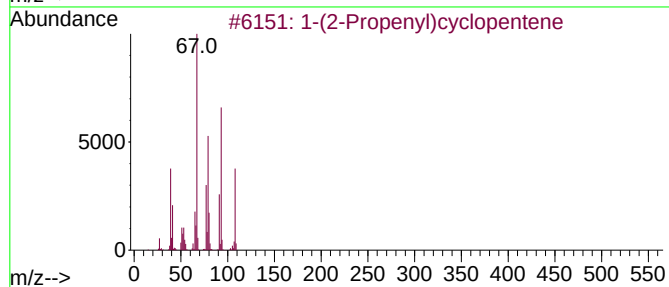
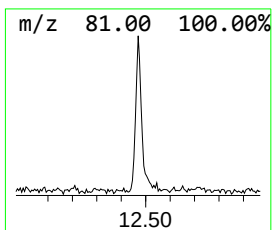
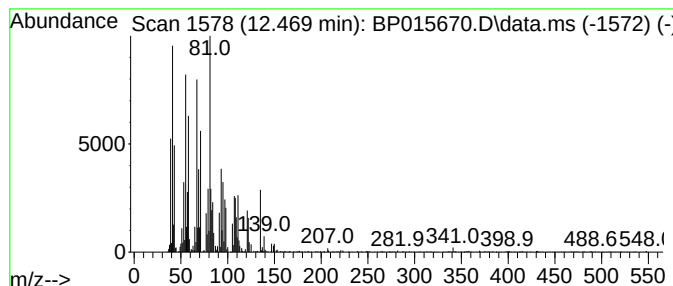
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.469	2.01 ng/ul	110365	Naphthalene-d8	10.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(2-Propenyl)cyclopentene	108	C8H12	037689-19-3	42
2			1-(2-Propenyl)cyclopentene	108	C8H12	037689-19-3	38
3			Undec-10-ynoic acid, undec-2-en-...	334	C22H38O2	1000406-96-8	35
4			4-Ethylcyclohexanol	128	C8H16O	004534-74-1	35
5			3-Oxatricyclo[3.2.1.0(2,4)]octan...	110	C7H10O	003146-39-2	25



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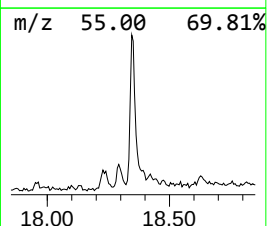
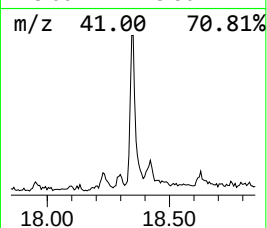
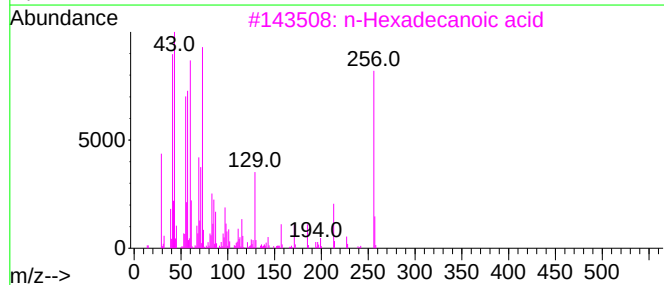
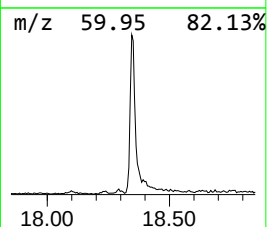
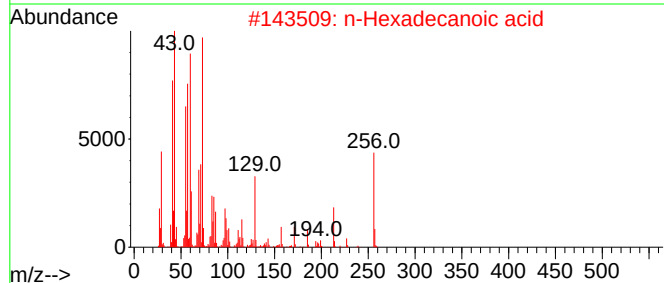
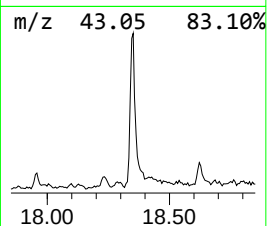
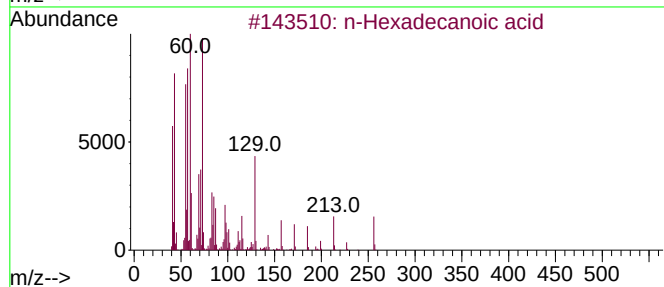
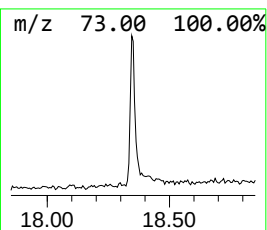
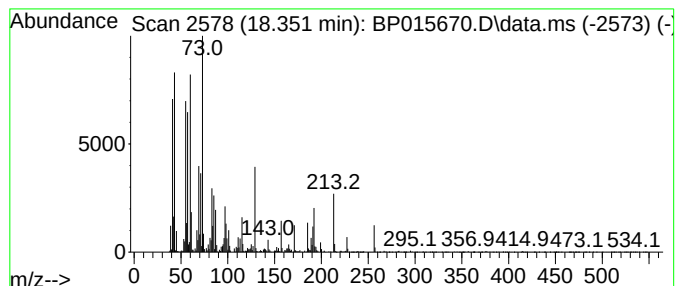
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.351	6.22 ng/ul	417597	Phenanthrene-d10	17.492

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	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015670.D
Acq On : 23 Jun 2023 14:23
Operator : MA/JU
Sample : 03084-19
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFN6

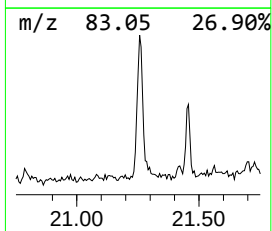
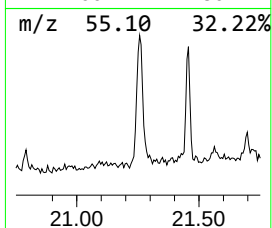
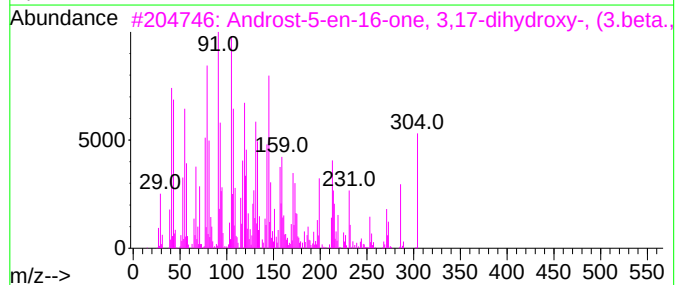
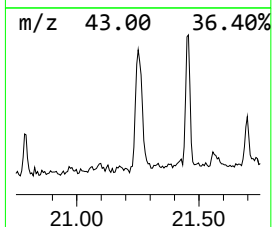
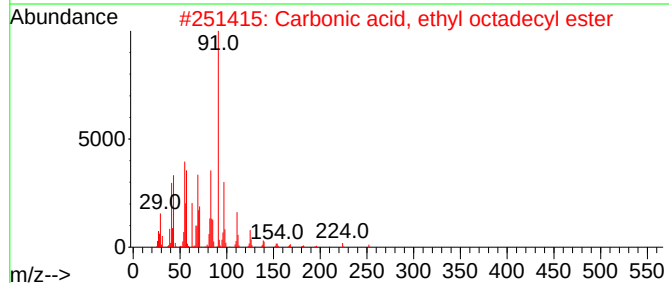
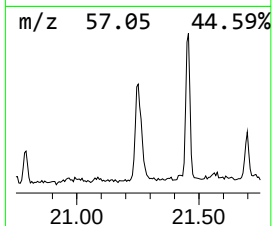
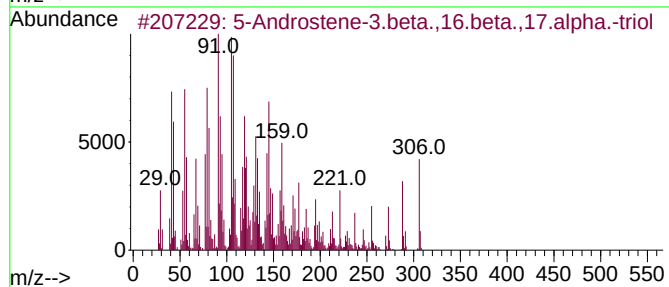
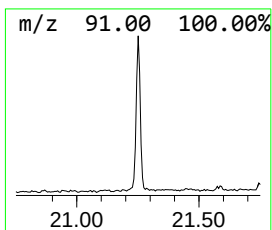
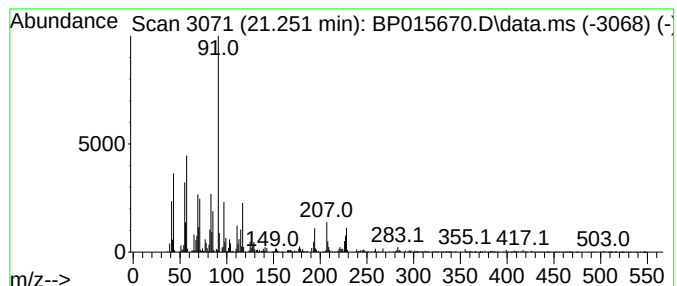
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	4.82 ng/ul	488279	Chrysene-d12	21.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Androstene-3.beta.,16.beta.,17...	306	C19H30O3	025126-70-9	22
2			Carbonic acid, ethyl octadecyl e...	342	C21H42O3	1000314-59-8	18
3			Androst-5-en-16-one, 3,17-dihydr...	304	C19H28O3	001159-66-6	15
4			Carbonic acid, ethyl tridecyl ester	272	C16H32O3	1000314-59-3	14
5			Cyclohexane, 1,3,5-triphenyl-	312	C24H24	028336-57-4	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015670.D
Acq On : 23 Jun 2023 14:23
Operator : MA/JU
Sample : 03084-19
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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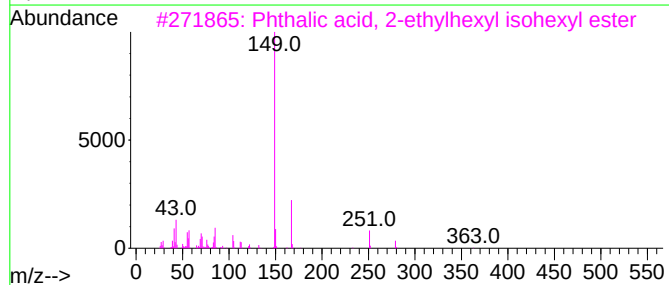
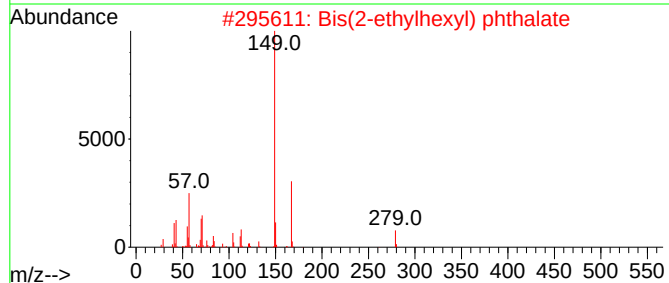
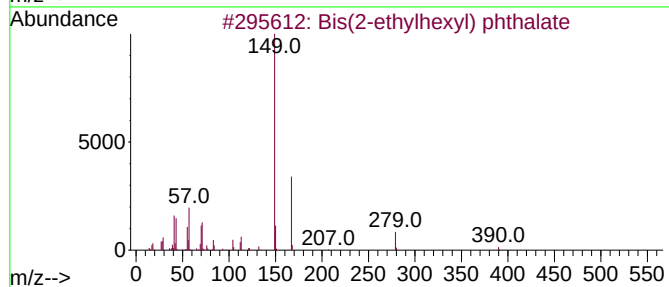
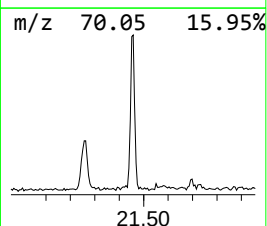
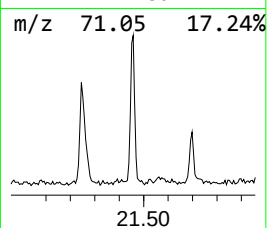
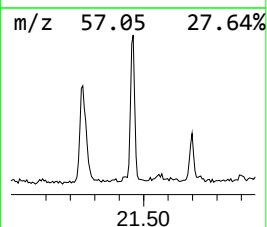
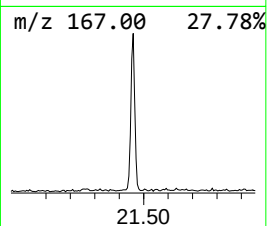
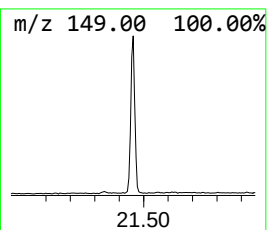
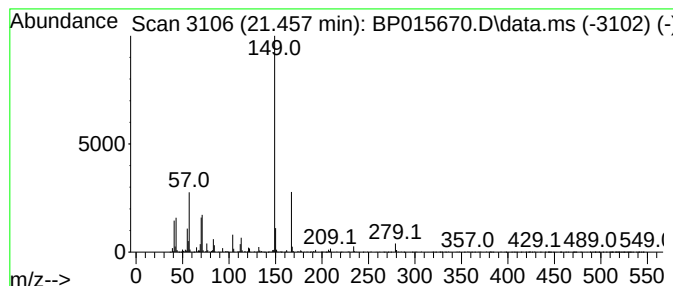
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Bis(2-ethylhexyl) phthalate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.457	5.22 ng/ul	528073	Chrysene-d12	21.586

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	86
3			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	86
4			Phthalic acid, di(oct-3-yl) ester	390	C24H38O4	1000377-72-3	86
5			Phthalic acid, di(hept-3-yl) ester	362	C22H34O4	1000357-00-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015670.D
Acq On : 23 Jun 2023 14:23
Operator : MA/JU
Sample : 03084-19
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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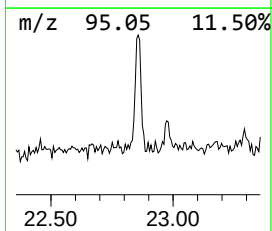
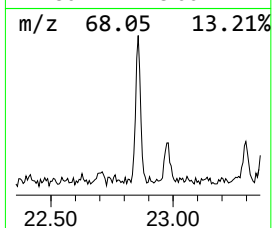
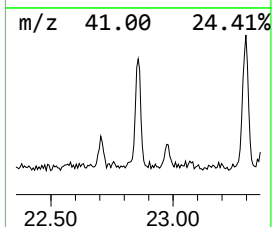
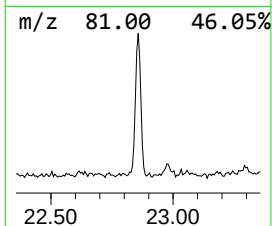
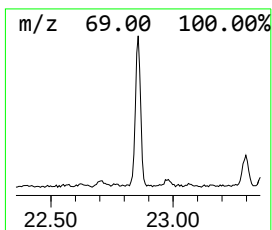
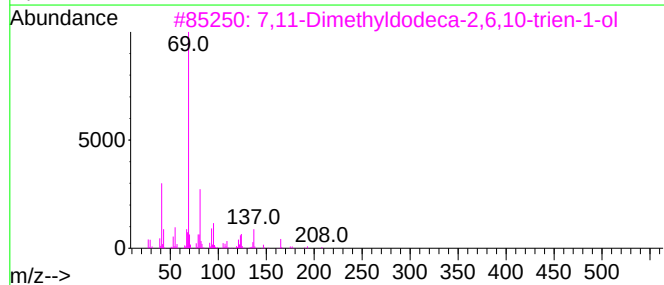
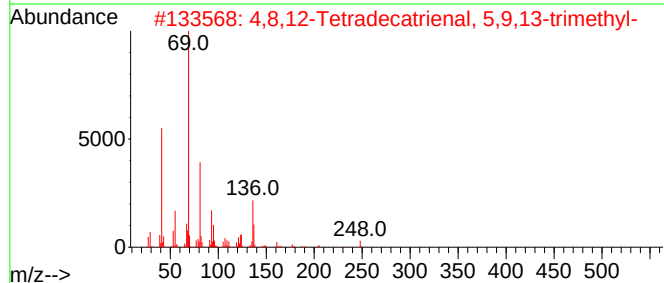
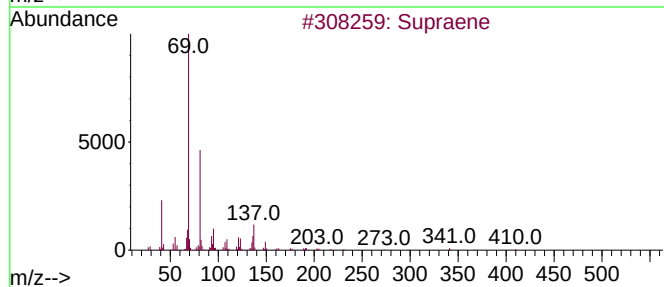
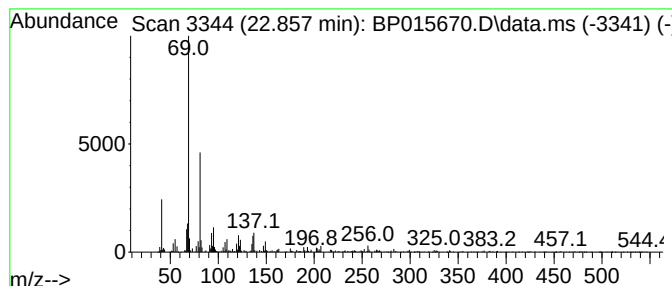
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.857	3.20 ng/ul	323760	Chrysene-d12	21.586

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	74
2			4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	64
3			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	64
4			1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	64
5			2,6,10,14-Hexadecatetraenoic aci...	332	C22H36O2	024035-35-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015670.D
Acq On : 23 Jun 2023 14:23
Operator : MA/JU
Sample : 03084-19
Misc :
ALS Vial : 4 Sample Multiplier: 1

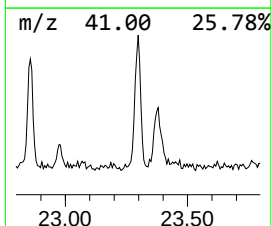
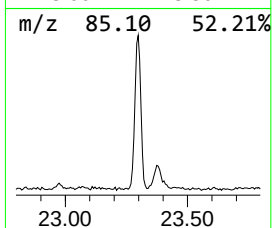
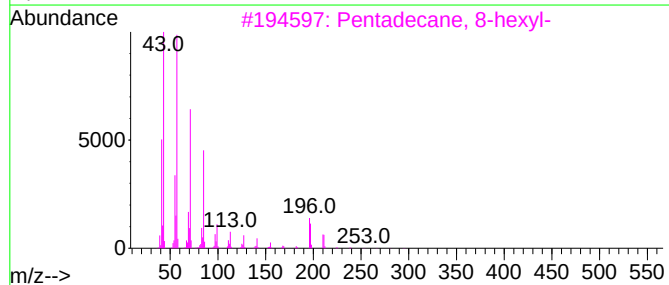
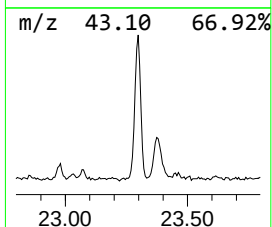
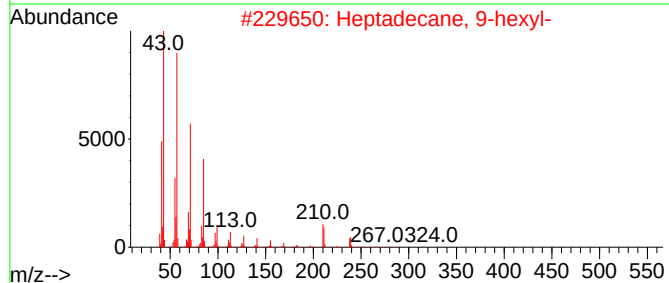
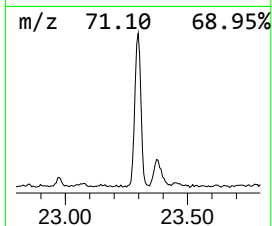
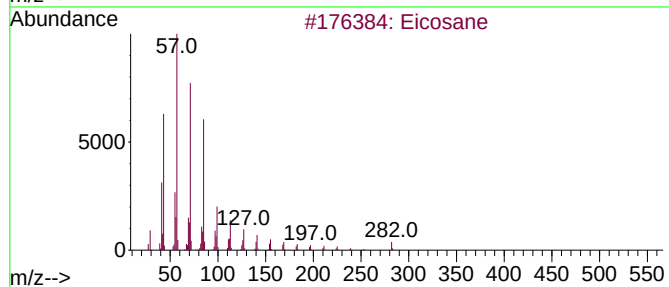
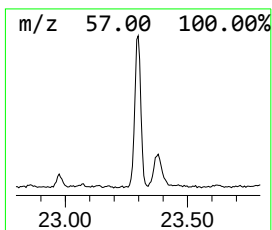
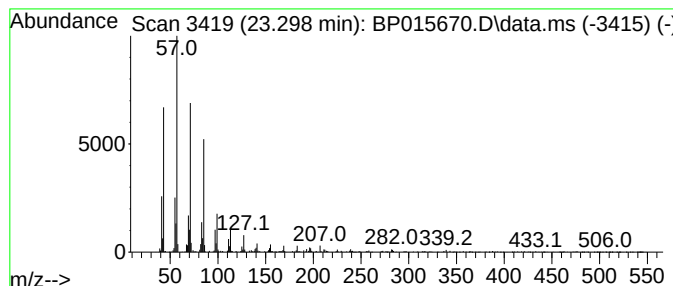
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
23.298	7.95 ng/ul	635306	Perylene-d12		24.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane		282	C20H42	000112-95-8	94
2	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	93
3	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	93
4	Hexacosane		366	C26H54	000630-01-3	90
5	Octacosane, 2-methyl-		408	C29H60	001560-98-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015670.D
Acq On : 23 Jun 2023 14:23
Operator : MA/JU
Sample : 03084-19
Misc :
ALS Vial : 4 Sample Multiplier: 1

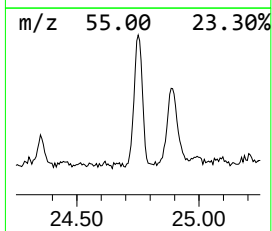
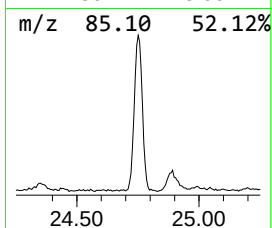
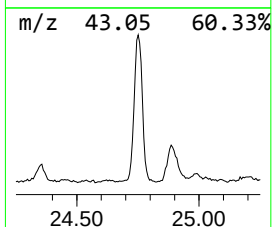
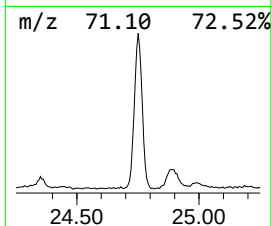
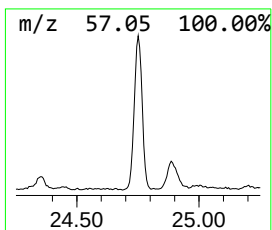
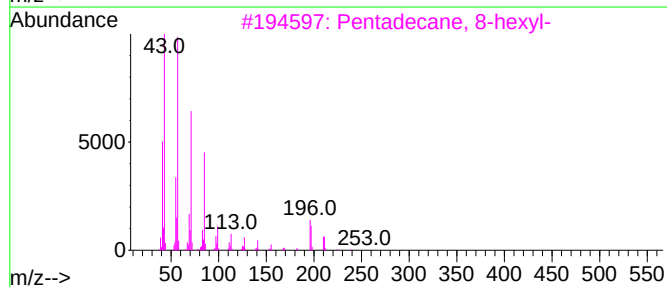
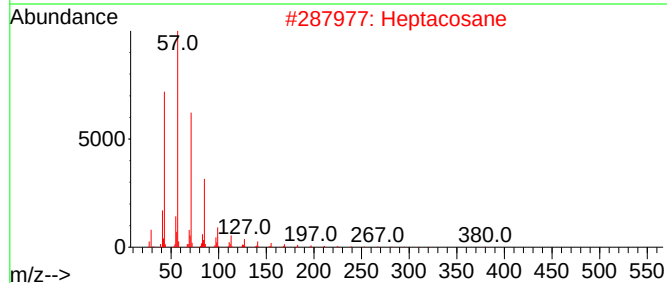
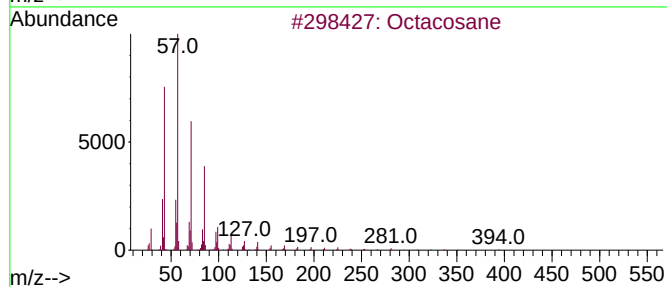
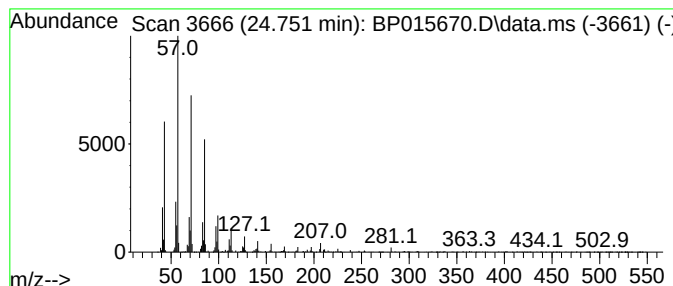
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.751	15.42 ng/ul	1231440	Perylene-d12		24.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	96
2	Heptacosane		380	C27H56	000593-49-7	93
3	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	93
4	Heneicosane, 11-pentyl-		366	C26H54	014739-72-1	93
5	2-Bromo dodecane		248	C12H25Br	013187-99-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062323\
Data File : BP015670.D
Acq On : 23 Jun 2023 14:23
Operator : MA/JU
Sample : 03084-19
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

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unknown-01	12.469	2.0	ng/u1	110365	2	10.910	1098570	20.0
n-Hexadecanoic ...	18.351	6.2	ng/u1	417597	4	17.492	1342680	20.0
unknown-02	21.251	4.8	ng/u1	488279	5	21.586	2024550	20.0
Bis(2-ethylhexy...	21.457	5.2	ng/u1	528073	5	21.586	2024550	20.0
Supraene	22.857	3.2	ng/u1	323760	5	21.586	2024550	20.0
(DEL) Alkane: S...	23.298	8.0	ng/u1	635306	6	24.145	1597590	20.0
(DEL) Alkane: S...	24.751	15.4	ng/u1	1231440	6	24.145	1597590	20.0