

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015560.D
 Acq On : 15 Jun 2023 17:35
 Operator : MA/JU
 Sample : 03088-01
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFQ9

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: BP015560.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	35	37	44	rVB3	10356	14895	0.46%	0.040%
2	3.958	127	131	138	rVB	27984	40236	1.25%	0.109%
3	4.081	148	152	159	rVB	14031	22116	0.69%	0.060%
4	4.181	165	169	172	rBV3	10835	13254	0.41%	0.036%
5	4.764	264	268	273	rVB8	4229	7028	0.22%	0.019%
6	5.128	327	330	338	rBV4	4332	7788	0.24%	0.021%
7	5.252	343	351	356	rBV10	6741	14363	0.45%	0.039%
8	6.046	481	486	493	rBV8	6169	14337	0.45%	0.039%
9	6.305	525	530	538	rBV	9075	19121	0.60%	0.052%
10	6.946	633	639	643	rBV8	2839	5951	0.19%	0.016%
11	7.099	662	665	667	rBV4	4014	5022	0.16%	0.014%
12	7.246	682	690	705	rBV2	195662	403087	12.57%	1.090%
13	7.410	711	718	727	rBV	162382	261220	8.14%	0.706%
14	7.610	745	752	763	rBV	205298	358405	11.17%	0.969%
15	7.687	763	765	769	rVB5	5085	5598	0.17%	0.015%
16	8.087	825	833	842	rBV	605186	985435	30.72%	2.665%
17	8.805	949	955	968	rBV	90854	182008	5.67%	0.492%
18	8.922	969	975	982	rVV	200895	333342	10.39%	0.901%
19	8.987	982	986	995	rVB	23793	52148	1.63%	0.141%
20	9.263	1027	1033	1046	rBV	225400	384828	12.00%	1.041%
21	9.422	1055	1060	1065	rVB9	4295	6671	0.21%	0.018%
22	9.640	1093	1097	1102	rVB8	4647	6374	0.20%	0.017%
23	9.993	1149	1157	1176	rBV	190841	352340	10.99%	0.953%
24	10.181	1184	1189	1197	rBV3	14340	43007	1.34%	0.116%
25	10.357	1214	1219	1228	rVB2	18189	34490	1.08%	0.093%
26	10.540	1243	1250	1267	rBV	227791	461781	14.40%	1.249%
27	10.910	1306	1313	1326	rVV	915686	1542150	48.08%	4.170%
28	11.063	1332	1339	1353	rBV	83360	199807	6.23%	0.540%
29	11.475	1403	1409	1411	rBV6	8801	15504	0.48%	0.042%
30	11.740	1448	1454	1462	rBV6	10058	28040	0.87%	0.076%
31	12.146	1518	1523	1531	rVB6	2394	5540	0.17%	0.015%
32	12.304	1543	1550	1554	rBV8	5029	9030	0.28%	0.024%
33	12.498	1579	1583	1589	rVB4	7288	12049	0.38%	0.033%
34	12.569	1589	1595	1599	rBV9	3793	5466	0.17%	0.015%
35	13.081	1677	1682	1687	rBV9	3963	7741	0.24%	0.021%
36	13.328	1719	1724	1731	rBV7	7034	9823	0.31%	0.027%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

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Max Peaks: 100

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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-BP060723.MA.M
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37	13.534	1755	1759	1764	rBV6	3309	4668	0.15%	0.013%
38	13.610	1768	1772	1778	rVB4	11530	20524	0.64%	0.056%
39	14.134	1854	1861	1868	rBV	601919	850228	26.51%	2.299%
40	14.245	1877	1880	1884	rVB6	3753	5071	0.16%	0.014%
41	14.428	1903	1911	1926	rVV	612470	954740	29.77%	2.582%
42	14.610	1935	1942	1946	rVB9	7018	12921	0.40%	0.035%
43	14.728	1952	1962	1971	rBV	1460179	2181784	68.03%	5.900%
44	14.963	1996	2002	2013	rBV2	57556	176414	5.50%	0.477%
45	15.134	2027	2031	2040	rVB2	12360	23849	0.74%	0.064%
46	15.510	2091	2095	2098	rBV5	5607	6909	0.22%	0.019%
47	15.551	2098	2102	2107	rVB6	7705	12805	0.40%	0.035%
48	15.722	2125	2131	2140	rBV	836019	1229712	38.34%	3.326%
49	15.857	2148	2154	2166	rBV	166463	287524	8.96%	0.778%
50	15.969	2169	2173	2180	rVB10	8782	16762	0.52%	0.045%
51	16.087	2186	2193	2202	rBV	95948	151472	4.72%	0.410%
52	16.422	2246	2250	2252	rBV3	13209	19712	0.61%	0.053%
53	16.575	2270	2276	2281	rBV9	9637	21503	0.67%	0.058%
54	16.651	2286	2289	2295	rVB6	9036	14308	0.45%	0.039%
55	16.869	2322	2326	2332	rBV9	9205	14244	0.44%	0.039%
56	16.945	2337	2339	2344	rVV4	5654	6465	0.20%	0.017%
57	17.016	2346	2351	2356	rVV3	15194	23459	0.73%	0.063%
58	17.216	2382	2385	2390	rVB4	9625	13918	0.43%	0.038%
59	17.275	2391	2395	2398	rVV6	3897	5853	0.18%	0.016%
60	17.363	2405	2410	2417	rBV10	21784	42254	1.32%	0.114%
61	17.428	2418	2421	2425	rVV5	11434	13681	0.43%	0.037%
62	17.486	2425	2431	2437	rVV	1908756	2718201	84.75%	7.351%
63	17.528	2437	2438	2442	rVV2	54081	63553	1.98%	0.172%
64	17.586	2442	2448	2461	rVB2	833323	1228258	38.30%	3.322%
65	17.757	2474	2477	2483	rVB	8911	14164	0.44%	0.038%
66	17.957	2507	2511	2513	rBV4	9898	13690	0.43%	0.037%
67	18.128	2538	2540	2544	rVB4	13510	13473	0.42%	0.036%
68	18.233	2554	2558	2563	rBV3	25324	38702	1.21%	0.105%
69	18.286	2563	2567	2572	rBV	120583	148350	4.63%	0.401%
70	18.345	2572	2577	2596	rVV	607743	883812	27.56%	2.390%
71	18.622	2621	2624	2631	rBV4	19164	33875	1.06%	0.092%
72	18.880	2666	2668	2670	rBV2	13009	12670	0.40%	0.034%
73	18.910	2670	2673	2679	rVB2	42360	63517	1.98%	0.172%
74	19.128	2706	2710	2714	rVB	50273	62874	1.96%	0.170%
75	19.175	2715	2718	2723	rVB2	30052	36297	1.13%	0.098%
76	19.228	2723	2727	2732	rBV3	34850	48422	1.51%	0.131%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	19.333	2740	2745	2747	rBV5	49264	87909	2.74%	0.238%
78	19.428	2759	2761	2763	rBV3	23474	23133	0.72%	0.063%
79	19.457	2763	2766	2767	rVV2	45032	49706	1.55%	0.134%
80	19.486	2767	2771	2778	rVB	236765	343640	10.71%	0.929%
81	19.569	2781	2785	2792	rBV2	171360	237754	7.41%	0.643%
82	19.810	2819	2826	2838	rVB2	1258896	1880824	58.64%	5.086%
83	20.298	2907	2909	2919	rVB5	64069	100920	3.15%	0.273%
84	20.616	2960	2963	2968	rBV7	45682	81389	2.54%	0.220%
85	20.780	2989	2991	2997	rVB	49412	60517	1.89%	0.164%
86	21.245	3066	3070	3076	rBV2	378645	602937	18.80%	1.631%
87	21.445	3102	3104	3108	rVB	96589	100132	3.12%	0.271%
88	21.569	3119	3125	3130	rBV	2314681	3207305	100.00%	8.674%
89	22.169	3224	3227	3230	rBV2	175661	217353	6.78%	0.588%
90	22.210	3230	3234	3241	rVB2	313897	580978	18.11%	1.571%
91	22.963	3359	3362	3367	rVB3	180937	262658	8.19%	0.710%
92	23.286	3412	3417	3422	rBV	632960	1019735	31.79%	2.758%
93	23.951	3524	3530	3537	rBV3	661135	1332468	41.54%	3.603%
94	24.115	3551	3558	3566	rBV	1335541	2820153	87.93%	7.627%
95	24.339	3591	3596	3604	rVB2	386190	770272	24.02%	2.083%
96	24.627	3640	3645	3652	rVB4	279798	593059	18.49%	1.604%
97	24.739	3658	3664	3674	rVB	586679	1245002	38.82%	3.367%
98	24.868	3680	3686	3701	rVB2	335290	1062504	33.13%	2.873%
99	26.198	3906	3912	3924	rVB2	546613	1472049	45.90%	3.981%
100	26.727	3996	4002	4022	rVB2	407623	1464844	45.67%	3.961%

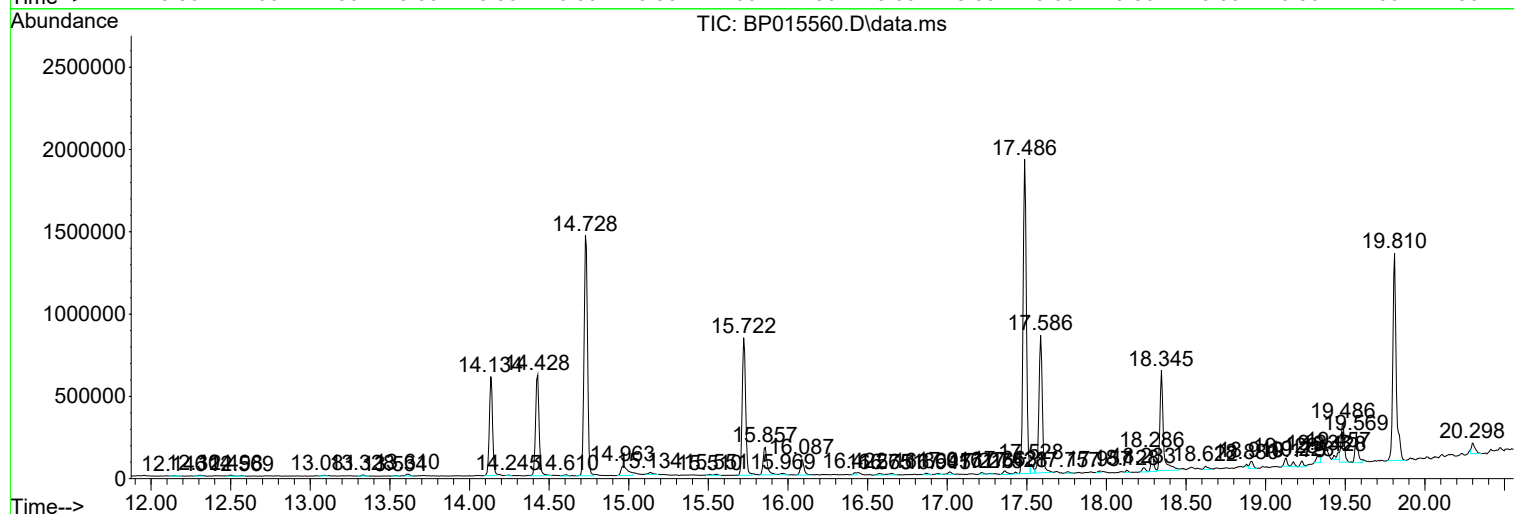
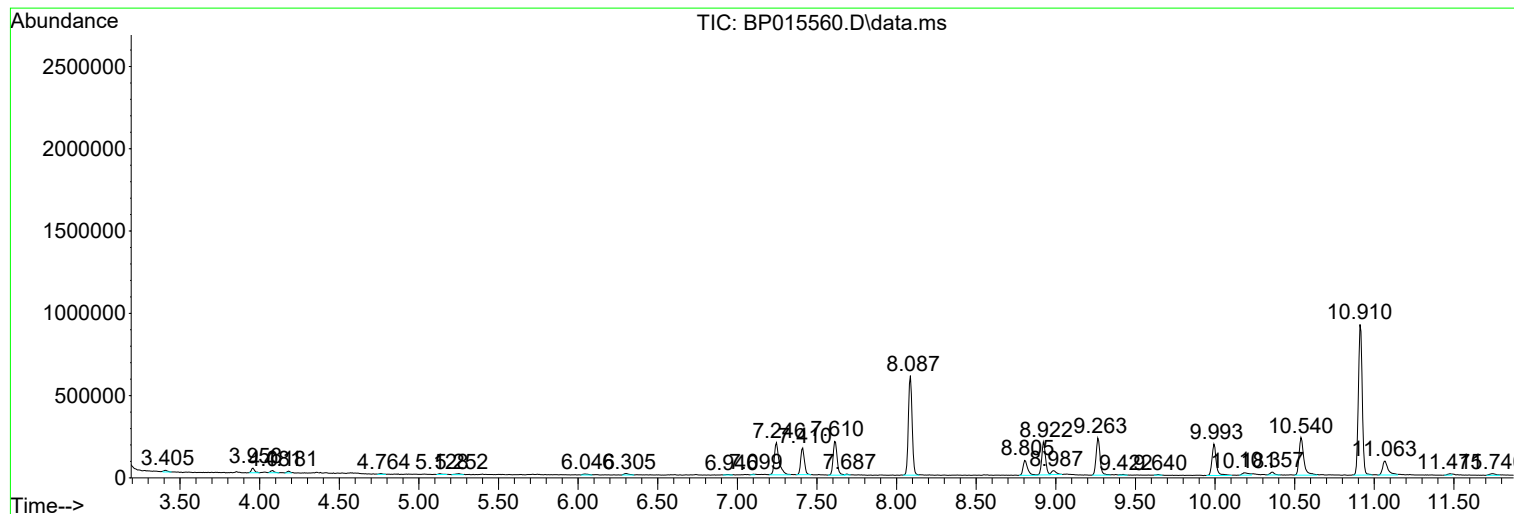
Sum of corrected areas: 36977874

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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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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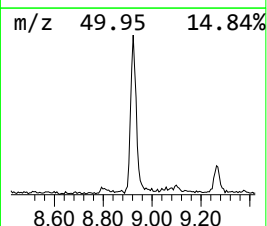
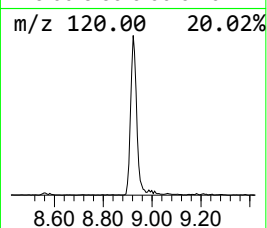
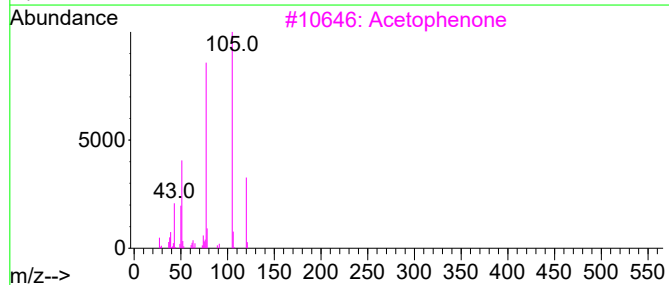
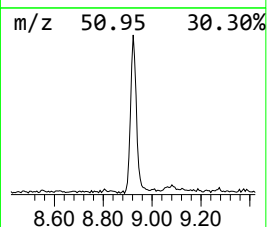
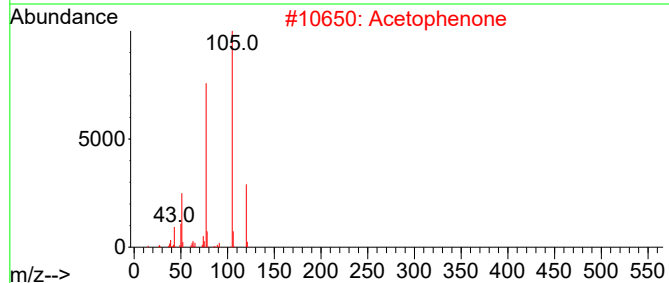
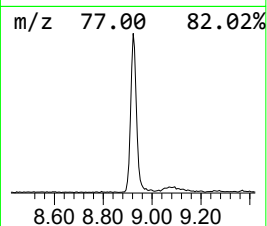
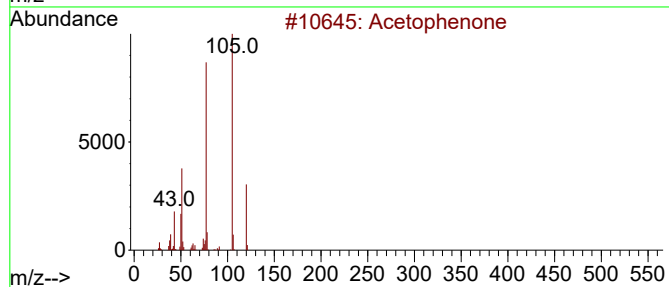
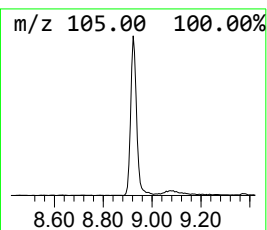
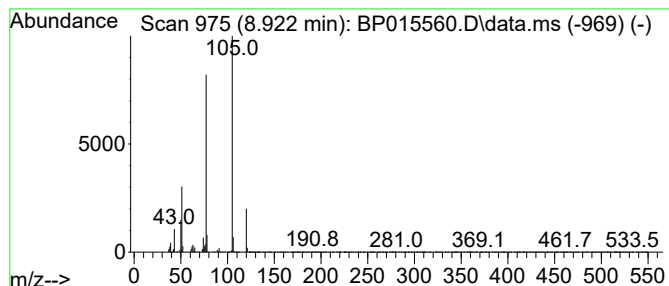
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	6.77 ng/ul	333342	1,4-Dichlorobenzene-d4	8.087

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	94
4	Acetophenone			120	C8H8O	000098-86-2	91
5	Acetophenone			120	C8H8O	000098-86-2	91



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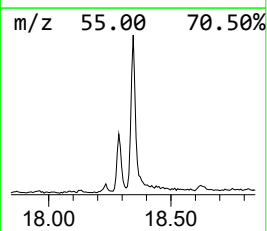
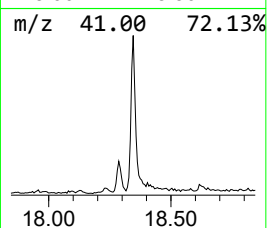
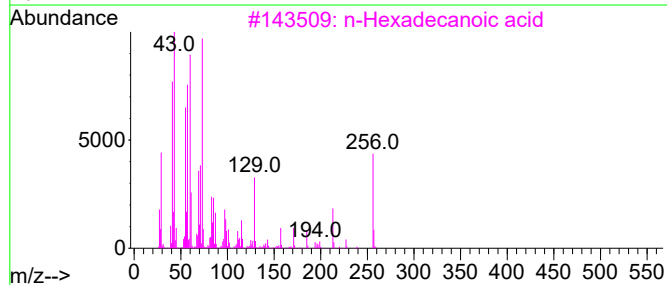
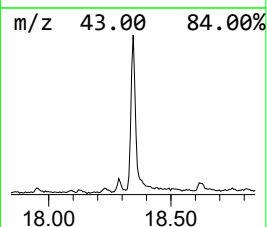
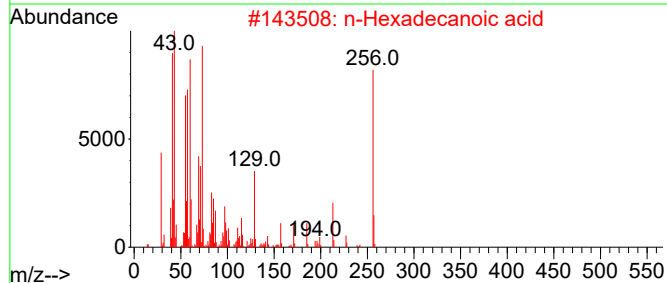
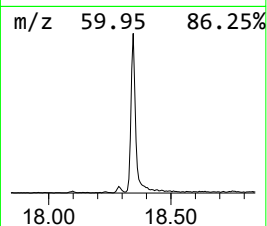
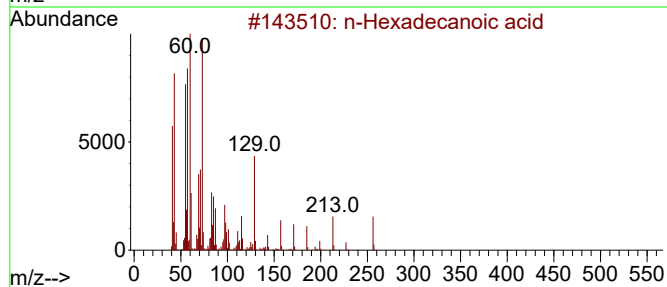
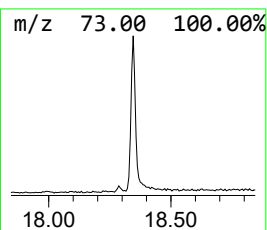
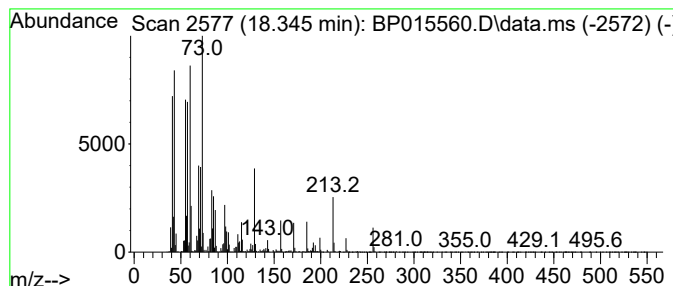
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	6.50 ng/ul	883812	Phenanthrene-d10	17.486

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	97
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	93



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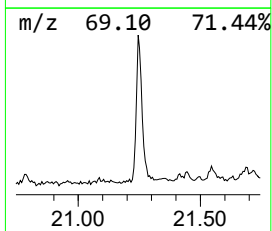
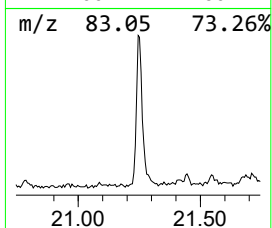
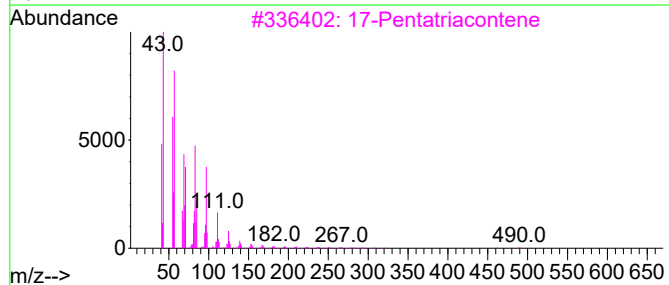
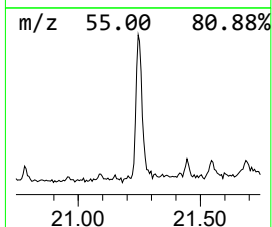
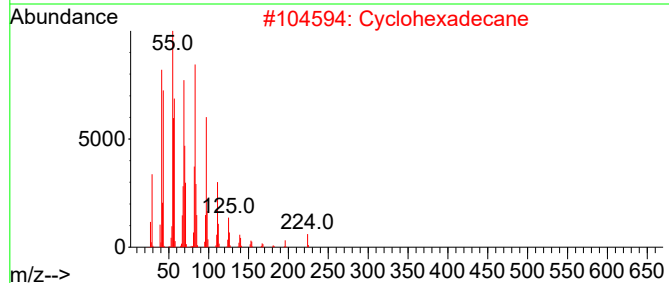
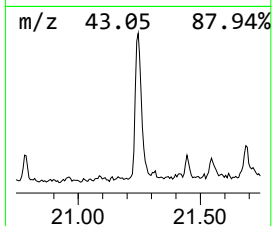
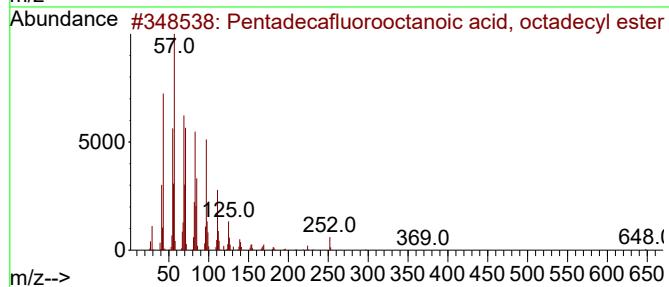
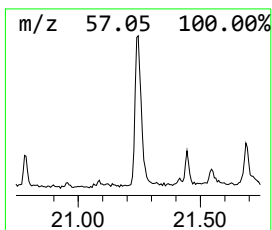
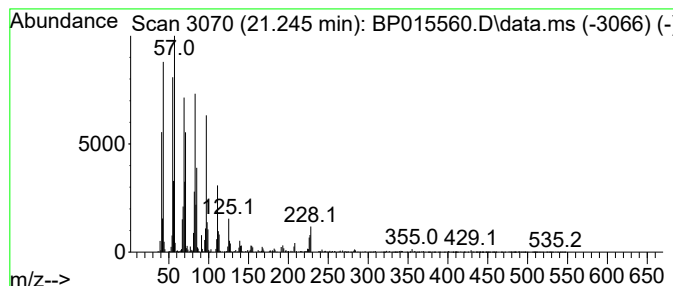
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentadecafluorooctanoic aci... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.245	3.76 ng/ul	602937	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94
2			Cyclohexadecane	224	C16H32	000295-65-8	92
3			17-Pentatriacontene	491	C35H70	006971-40-0	91
4			1-Hexacosene	364	C26H52	018835-33-1	91
5			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
Acq On : 15 Jun 2023 17:35
Operator : MA/JU
Sample : 03088-01
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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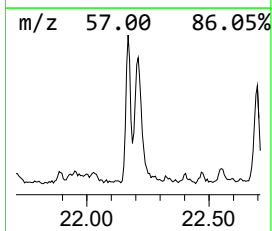
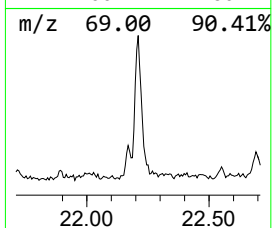
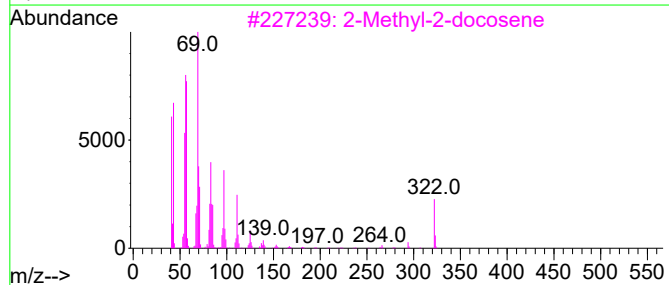
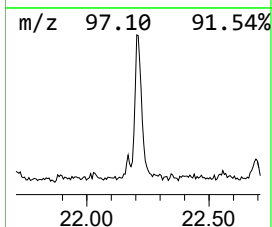
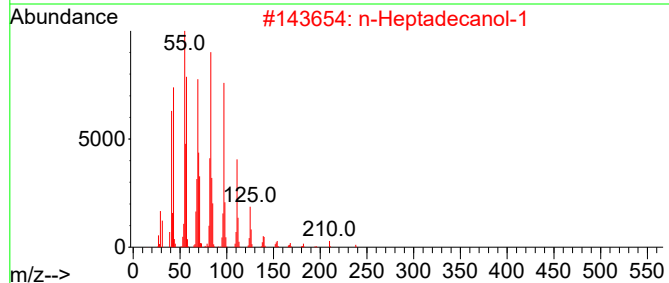
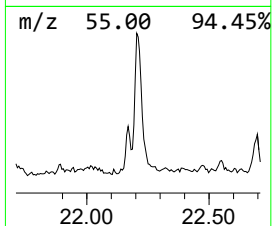
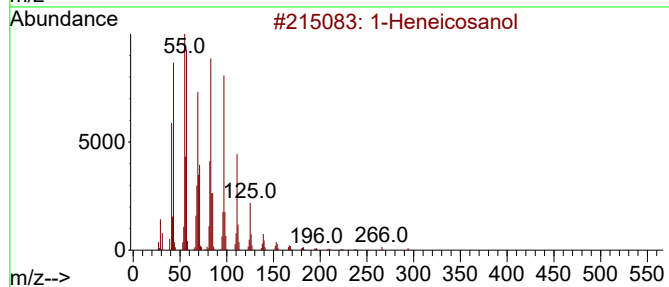
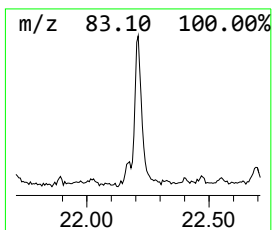
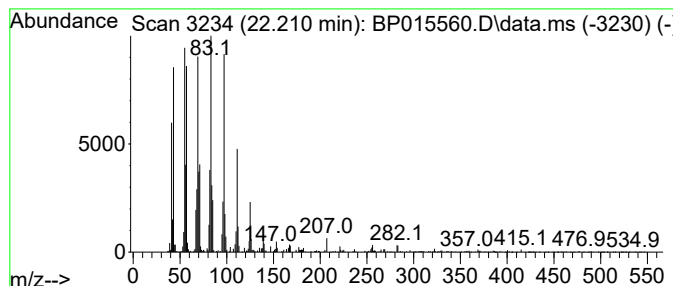
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Heneicosanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.210	3.62 ng/ul	580978	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol			312	C21H44O	015594-90-8	95
2	n-Heptadecanol-1			256	C17H36O	001454-85-9	94
3	2-Methyl-2-docosene			322	C23H46	1000131-16-9	93
4	n-Nonadecanol-1			284	C19H40O	001454-84-8	91
5	Octacosanol			410	C28H58O	000557-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
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Sample : 03088-01
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Instrument :
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ClientSampleId :
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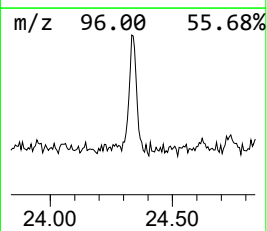
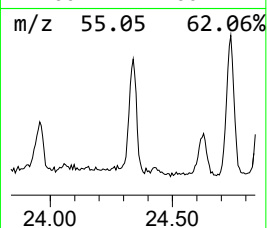
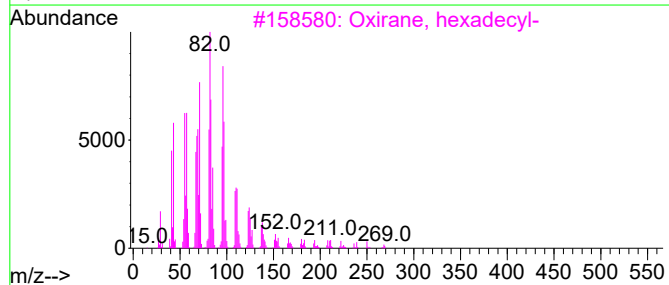
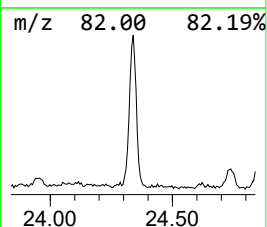
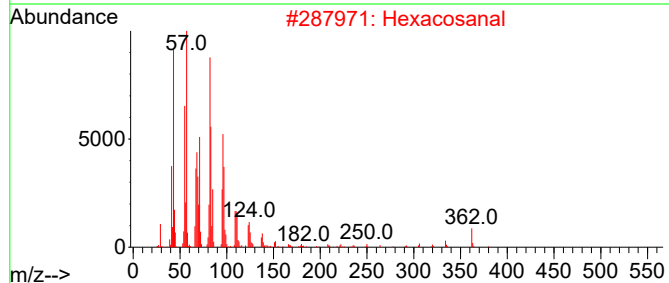
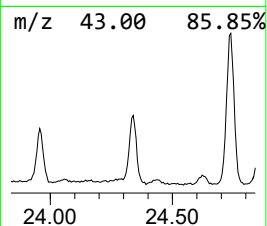
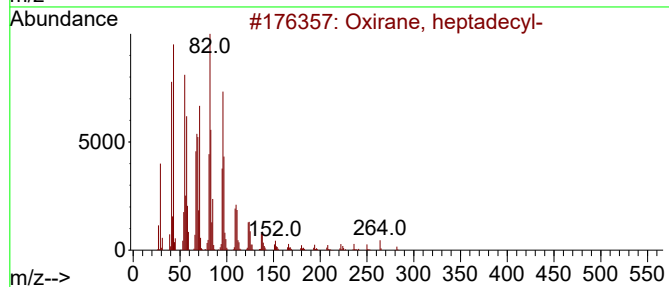
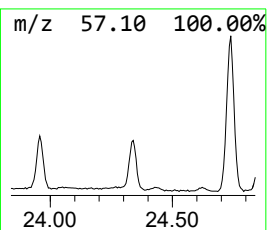
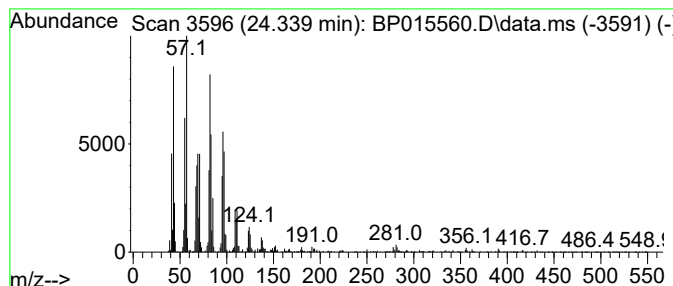
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Oxirane, heptadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.339	5.46 ng/ul	770272	Perylene-d12	24.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94
2			Hexacosanal	380	C26H52O	026627-85-0	91
3			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
4			Octadecanal	268	C18H36O	000638-66-4	91
5			(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
Acq On : 15 Jun 2023 17:35
Operator : MA/JU
Sample : 03088-01
Misc :
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Instrument :
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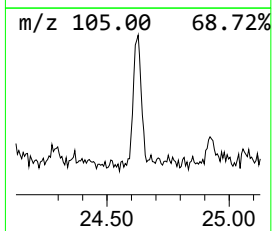
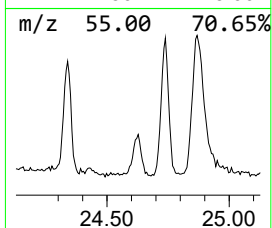
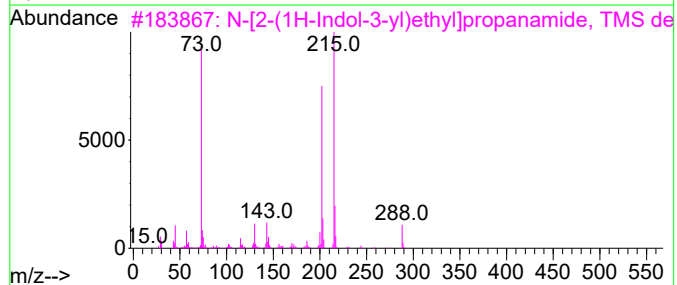
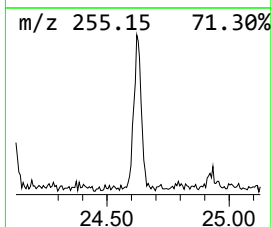
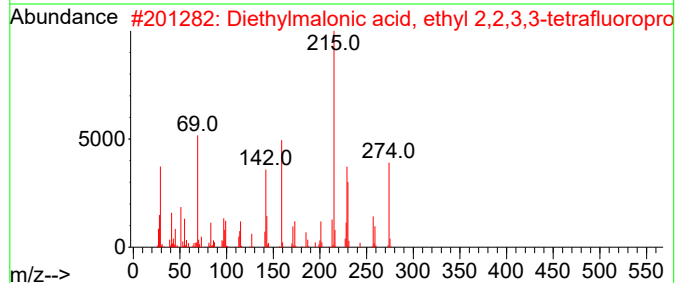
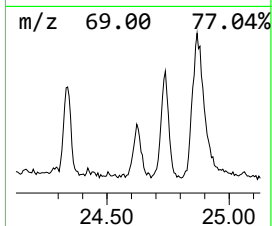
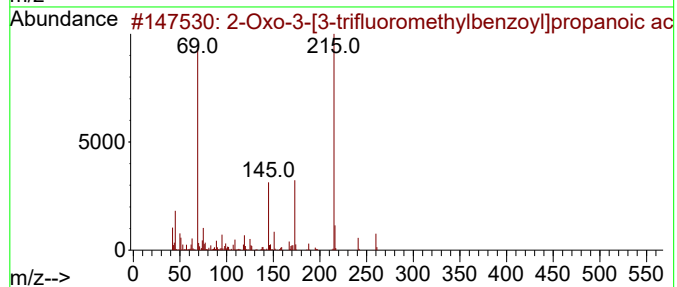
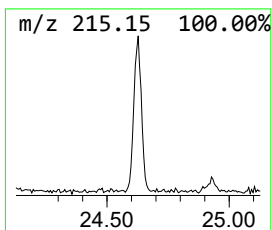
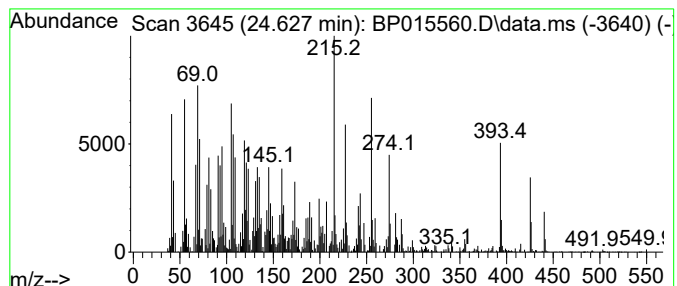
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.627	4.21 ng/ul	593059	Perylene-d12	24.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Oxo-3-[3-trifluoromethylbenzoyl]propanoic acid	260	C11H7F3O4	1000211-47-7	35
2			Diethylmalonic acid, ethyl 2,2,3,3-tetrafluoropropanoate	302	C12H18F4O4	1000370-82-9	27
3			N-[2-(1H-Indol-3-yl)ethyl]propanamide, TMS derivative	288	C16H24N2OSi	1000493-19-0	25
4			Ethanone, 1-[2,3-dihydro-1,1,2,6-tetrafluorocyclohex-5-en-1-yl]	258	C18H26O	068140-48-7	25
5			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
Acq On : 15 Jun 2023 17:35
Operator : MA/JU
Sample : 03088-01
Misc :
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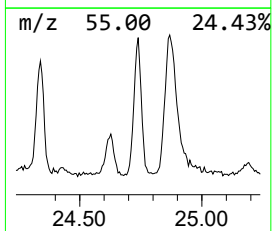
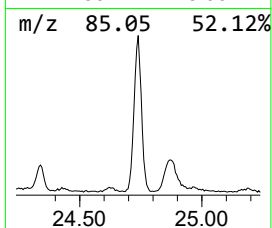
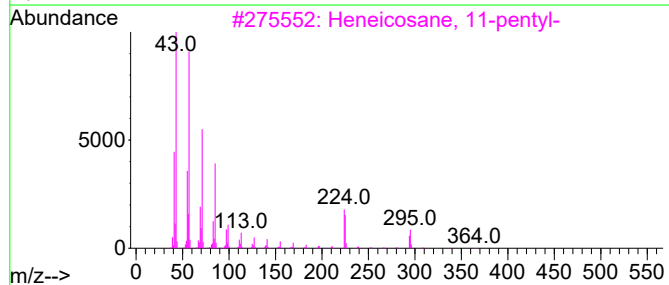
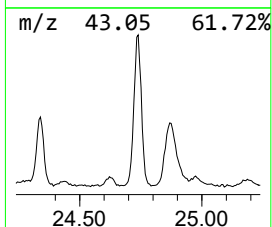
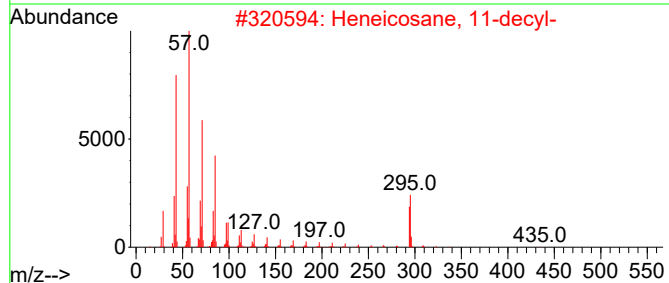
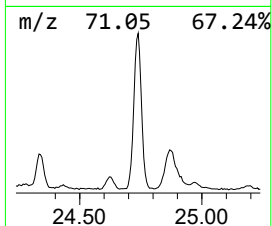
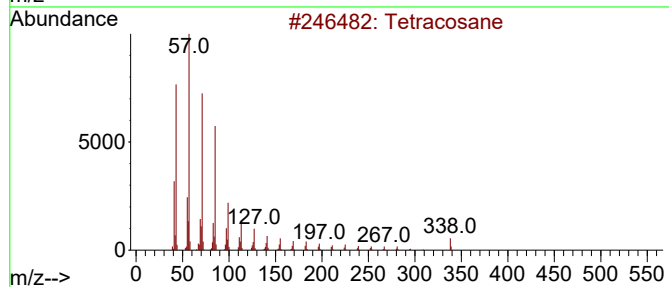
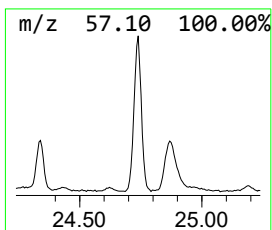
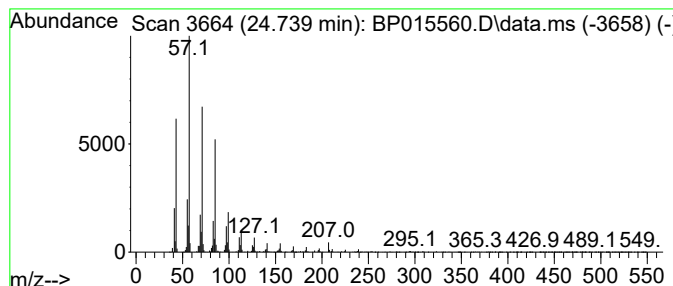
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.739	8.83 ng/ul	1245000	Perylene-d12	24.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	97
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	96
3		Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
5		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
Acq On : 15 Jun 2023 17:35
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Sample : 03088-01
Misc :
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Instrument :
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ClientSampleId :
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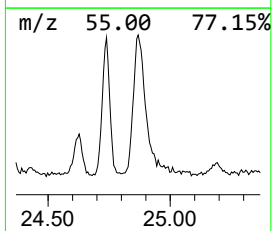
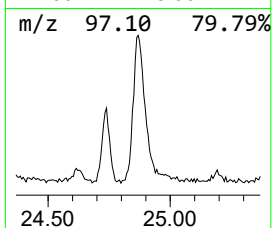
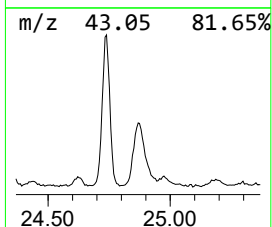
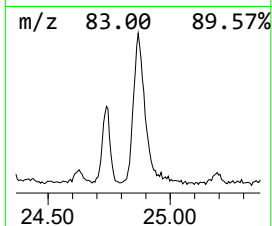
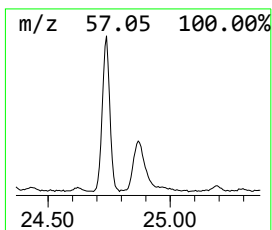
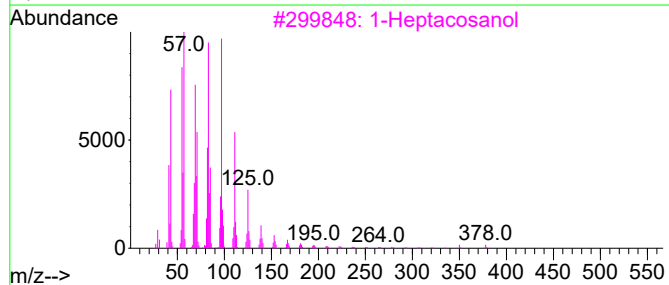
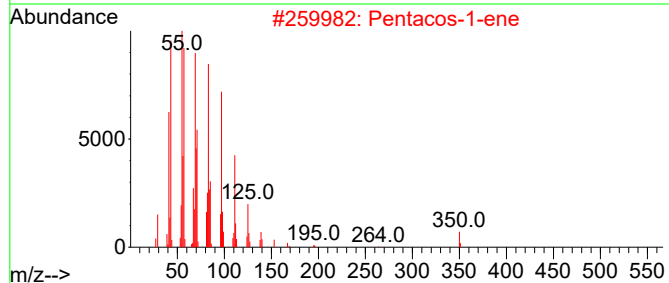
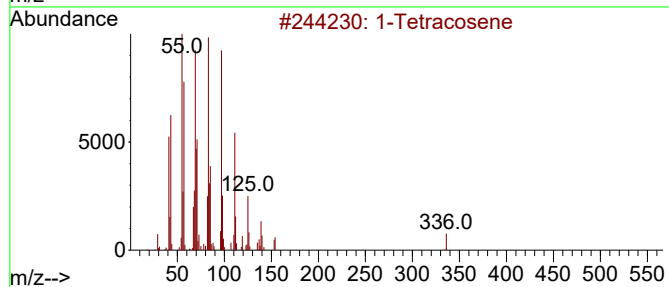
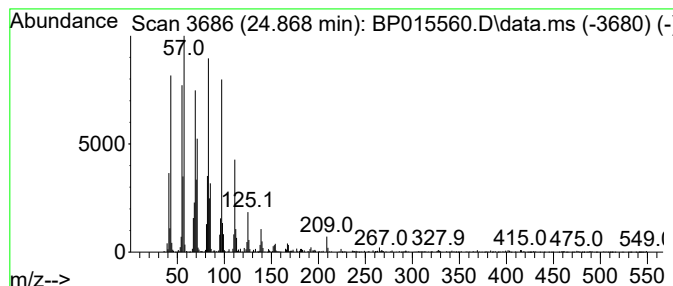
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.868	7.54 ng/ul	1062500	Perylene-d12	24.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tetracosene	336	C24H48	010192-32-2	96
2		Pentacos-1-ene	350	C25H50	016980-85-1	96
3		1-Heptacosanol	396	C27H56O	002004-39-9	94
4		1-Heneicosanol	312	C21H44O	015594-90-8	93
5		n-Tetracosanol-1	354	C24H50O	000506-51-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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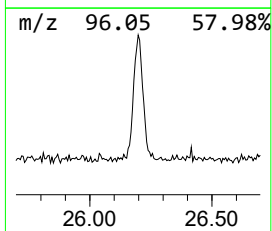
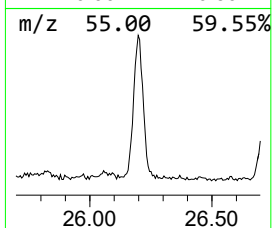
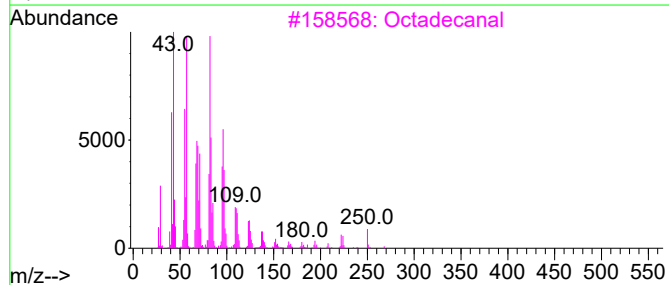
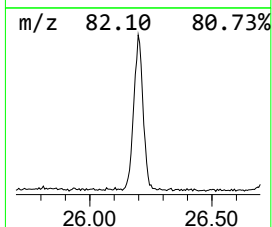
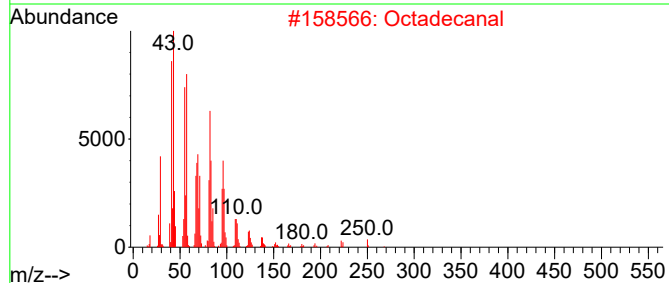
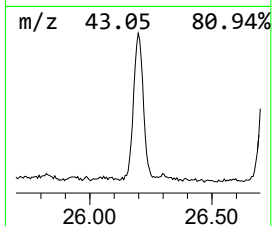
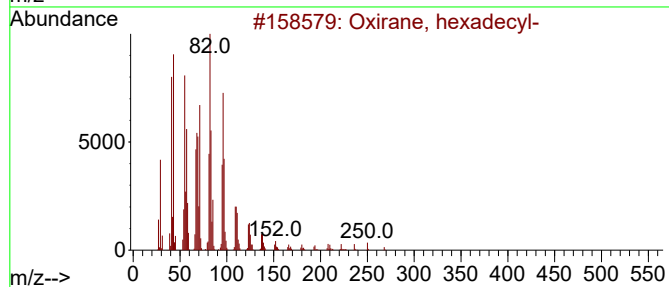
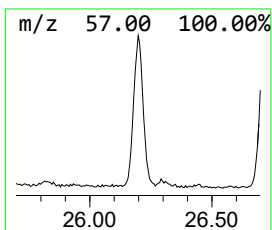
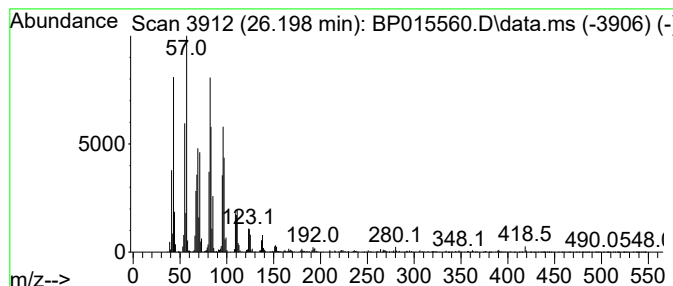
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Oxirane, hexadecyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.198	10.44 ng/ul	1472050	Perylene-d12	24.116	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
2	Octadecanal	268	C18H36O	000638-66-4	94
3	Octadecanal	268	C18H36O	000638-66-4	94
4	Dotriacontanal	464	C32H64O	057878-00-9	91
5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
Acq On : 15 Jun 2023 17:35
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ALS Vial : 54 Sample Multiplier: 1

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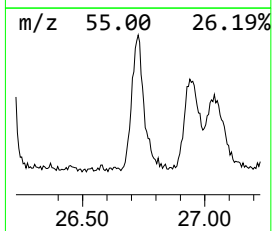
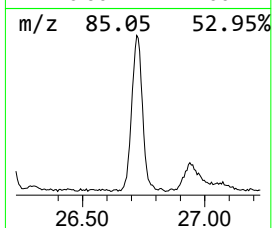
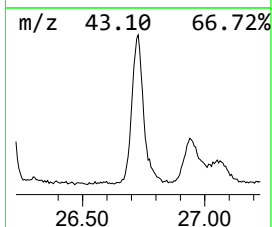
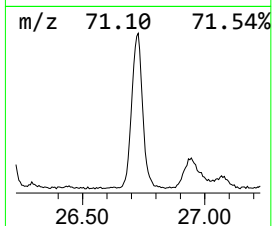
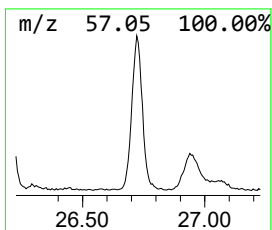
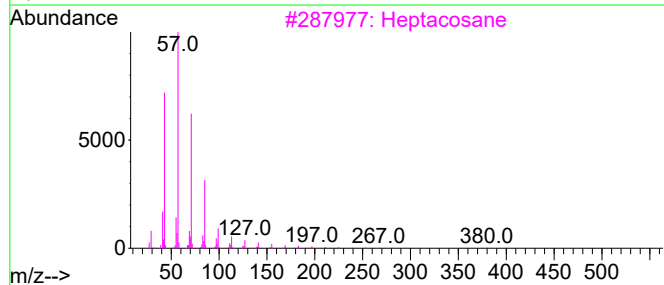
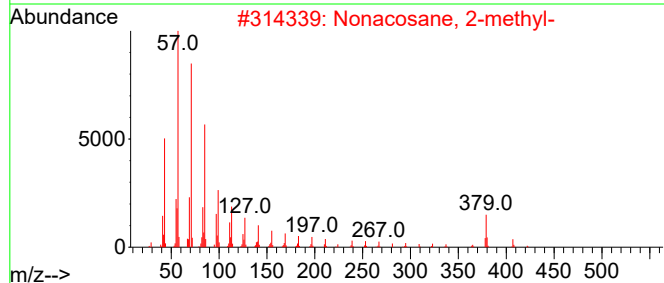
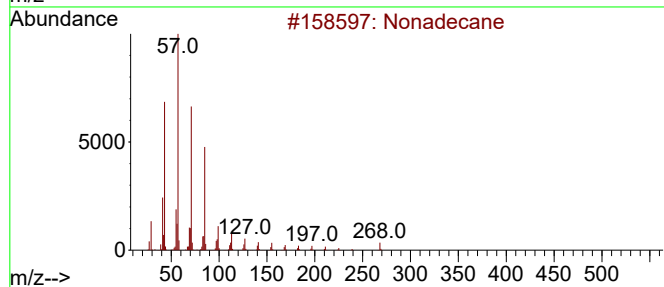
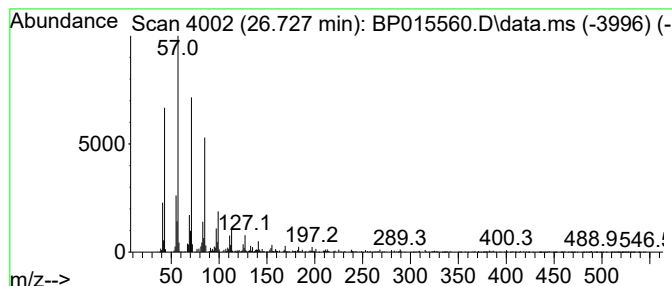
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.727	10.39 ng/ul	1464840	Perylene-d12	24.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	94
2		Nonacosane, 2-methyl-	422	C30H62	001560-75-4	93
3		Heptacosane	380	C27H56	000593-49-7	93
4		Tetracosane	338	C24H50	000646-31-1	93
5		Docosane, 11-butyl-	366	C26H54	013475-76-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015560.D
Acq On : 15 Jun 2023 17:35
Operator : MA/JU
Sample : 03088-01
Misc :
ALS Vial : 54 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFQ9

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Aceto phenone	8.922	6.8	ng/ul	333342	1	8.087	985435	20.0
n-Hexadecanoic ...	18.345	6.5	ng/ul	883812	4	17.486	2718200	20.0
Pentadecafluoro...	21.245	3.8	ng/ul	602937	5	21.569	3207310	20.0
1-Heneicosanol	22.210	3.6	ng/ul	580978	5	21.569	3207310	20.0
Oxirane, heptad...	24.339	5.5	ng/ul	770272	6	24.116	2820150	20.0
unknown-01	24.627	4.2	ng/ul	593059	6	24.116	2820150	20.0
(DEL) Alkane: S...	24.739	8.8	ng/ul	1245000	6	24.116	2820150	20.0
(DEL) Alkane: S...	24.868	7.5	ng/ul	1062500	6	24.116	2820150	20.0
Oxirane, hexade...	26.198	10.4	ng/ul	1472050	6	24.116	2820150	20.0
(DEL) Alkane: S...	26.727	10.4	ng/ul	1464840	6	24.116	2820150	20.0