

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
 Data File : BP015548.D
 Acq On : 15 Jun 2023 08:53
 Operator : MA/JU
 Sample : 03088-07
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFR3

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: BP015548.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	42	rVB3	8698	12487	0.34%	0.043%
2	3.916	121	124	125	rBV3	5756	7032	0.19%	0.024%
3	4.081	148	152	162	rVB2	13389	28229	0.77%	0.096%
4	4.416	206	209	213	rBV4	8307	10385	0.28%	0.035%
5	5.010	307	310	313	rVB5	3594	4650	0.13%	0.016%
6	5.134	326	331	334	rBV5	4284	7270	0.20%	0.025%
7	6.046	482	486	494	rVV10	4806	9646	0.26%	0.033%
8	6.304	523	530	536	rVV6	7040	17474	0.48%	0.060%
9	6.746	599	605	610	rVB5	8752	14594	0.40%	0.050%
10	7.246	682	690	702	rBV	145728	305036	8.31%	1.041%
11	7.410	712	718	730	rVV	119613	199179	5.42%	0.680%
12	7.510	731	735	740	rVB8	3850	6867	0.19%	0.023%
13	7.616	746	753	761	rBV	152552	271606	7.40%	0.927%
14	8.087	826	833	844	rBV	312547	520381	14.17%	1.777%
15	8.804	947	955	968	rBV	163610	311584	8.48%	1.064%
16	8.928	968	976	982	rBV	125223	223400	6.08%	0.763%
17	9.204	1019	1023	1025	rVB5	3176	3703	0.10%	0.013%
18	9.263	1027	1033	1047	rVV	149403	279574	7.61%	0.954%
19	9.428	1056	1061	1062	rBV5	3980	5981	0.16%	0.020%
20	9.992	1148	1157	1168	rBV	129533	249715	6.80%	0.853%
21	10.210	1187	1194	1204	rBV	10580	41843	1.14%	0.143%
22	10.357	1215	1219	1228	rVB3	12073	23819	0.65%	0.081%
23	10.539	1243	1250	1274	rBV	160671	362031	9.86%	1.236%
24	10.910	1306	1313	1326	rVB	486423	844068	22.98%	2.882%
25	11.069	1332	1340	1354	rBV	82115	197010	5.36%	0.673%
26	11.381	1391	1393	1398	rVB5	2510	3768	0.10%	0.013%
27	11.475	1401	1409	1417	rBV5	6306	19953	0.54%	0.068%
28	11.751	1449	1456	1462	rBV9	6341	18425	0.50%	0.063%
29	12.139	1519	1522	1528	rBV8	3534	5919	0.16%	0.020%
30	12.304	1547	1550	1553	rBV5	2653	3831	0.10%	0.013%
31	12.498	1579	1583	1585	rVV5	4139	6647	0.18%	0.023%
32	13.045	1673	1676	1679	rBV5	2461	3746	0.10%	0.013%
33	13.081	1679	1682	1691	rVB5	6974	14244	0.39%	0.049%
34	13.333	1718	1725	1730	rBV8	6011	9653	0.26%	0.033%
35	13.498	1748	1753	1755	rBV6	2336	4137	0.11%	0.014%
36	13.610	1763	1772	1780	rBV2	35502	59603	1.62%	0.203%

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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
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37	14.045	1840	1846	1851	rBV10	9444	14123	0.38%	0.048%
38	14.133	1853	1861	1869	rBV	414141	610861	16.63%	2.086%
39	14.245	1877	1880	1884	rVB6	4531	6241	0.17%	0.021%
40	14.369	1897	1901	1902	rBV4	3891	5054	0.14%	0.017%

41	14.422	1904	1910	1925	rVV	438128	700208	19.06%	2.391%
42	14.563	1930	1934	1938	rVV6	7640	12221	0.33%	0.042%
43	14.604	1938	1941	1944	rVB4	3698	4541	0.12%	0.016%
44	14.675	1949	1953	1956	rBV3	14519	18988	0.52%	0.065%
45	14.728	1956	1962	1969	rBV	823453	1205727	32.83%	4.116%

46	14.969	1994	2003	2021	rBV8	41442	161158	4.39%	0.550%
47	15.139	2029	2032	2037	rBV6	9547	14422	0.39%	0.049%
48	15.633	2110	2116	2120	rBV7	13651	21967	0.60%	0.075%
49	15.722	2124	2131	2143	rBV	587456	917607	24.98%	3.133%
50	15.857	2148	2154	2166	rVV	105791	208692	5.68%	0.712%

51	15.963	2169	2172	2176	rVB6	8387	12061	0.33%	0.041%
52	16.086	2188	2193	2200	rBV	75572	116180	3.16%	0.397%
53	16.286	2224	2227	2233	rVB7	8670	12348	0.34%	0.042%
54	16.445	2247	2254	2259	rBV7	12227	27847	0.76%	0.095%
55	16.580	2273	2277	2280	rBV6	9887	14378	0.39%	0.049%

56	17.269	2390	2394	2398	rBV6	4839	9833	0.27%	0.034%
57	17.363	2406	2410	2411	rBV4	14523	17119	0.47%	0.058%
58	17.427	2418	2421	2425	rBV6	10782	15749	0.43%	0.054%
59	17.486	2425	2431	2441	rVV	1064921	1615744	43.99%	5.516%
60	17.586	2443	2448	2461	rVB2	607668	907133	24.70%	3.097%

61	17.821	2485	2488	2493	rBV7	9600	15032	0.41%	0.051%
62	18.127	2537	2540	2544	rVB6	11001	12070	0.33%	0.041%
63	18.233	2555	2558	2562	rVB4	19972	24808	0.68%	0.085%
64	18.286	2563	2567	2572	rBV	84830	114192	3.11%	0.390%
65	18.345	2572	2577	2586	rBV	478472	656348	17.87%	2.241%

66	19.121	2707	2709	2715	rVB	45271	57429	1.56%	0.196%
67	19.427	2759	2761	2763	rBV3	31314	32610	0.89%	0.111%
68	19.457	2763	2766	2768	rVV3	69803	87278	2.38%	0.298%
69	19.480	2768	2770	2778	rVB	104616	150692	4.10%	0.514%
70	19.568	2781	2785	2796	rBV	180927	253783	6.91%	0.866%

71	19.810	2820	2826	2836	rBV	685799	1014692	27.63%	3.464%
72	21.245	3066	3070	3080	rBV4	251284	433323	11.80%	1.479%
73	21.568	3119	3125	3131	rBV	1285945	1831076	49.86%	6.251%
74	22.080	3201	3212	3224	rVB5	510037	2252444	61.33%	7.690%
75	23.286	3412	3417	3423	rVB	777351	1242374	33.83%	4.242%

76	23.951	3524	3530	3539	rVB2	455943	1005398	27.37%	3.432%
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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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Title : SVOA CALIBRATION

77	24.115	3552	3558	3572	rVB	738067	1641559	44.70%	5.604%
78	24.627	3639	3645	3653	rVB3	439225	926597	25.23%	3.163%
79	24.739	3657	3664	3675	rVB	1644746	3672749	100.00%	12.539%
80	24.862	3679	3685	3692	rBV5	178896	562755	15.32%	1.921%
81	26.727	3994	4002	4026	rVB	729795	2545652	69.31%	8.691%

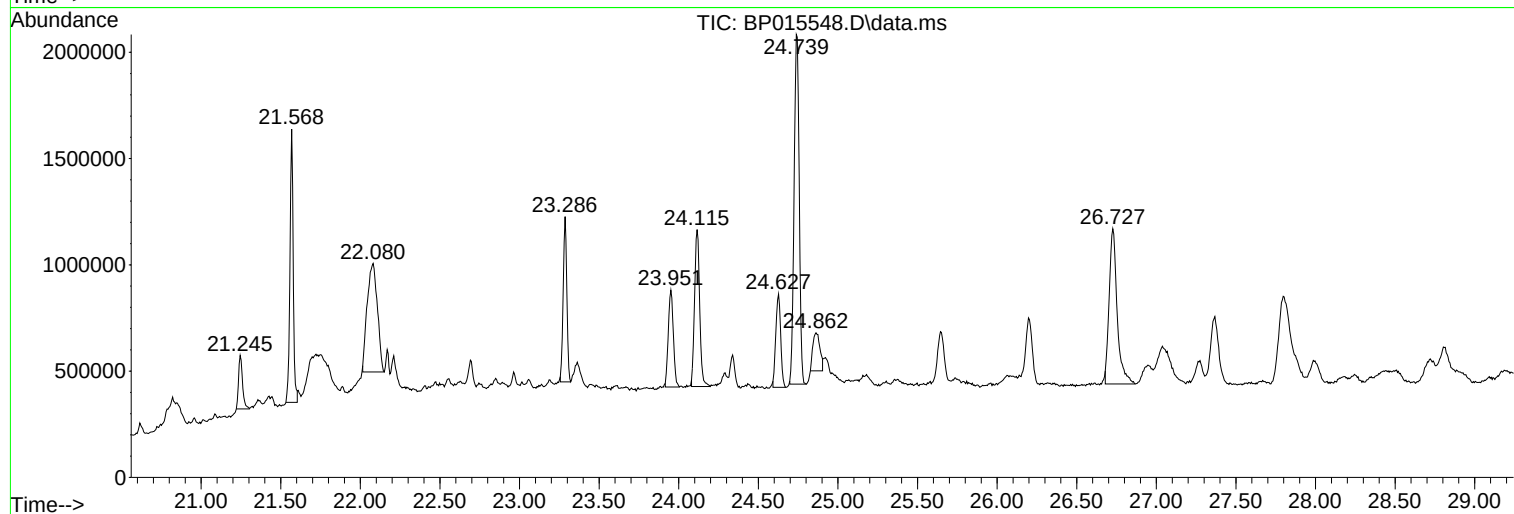
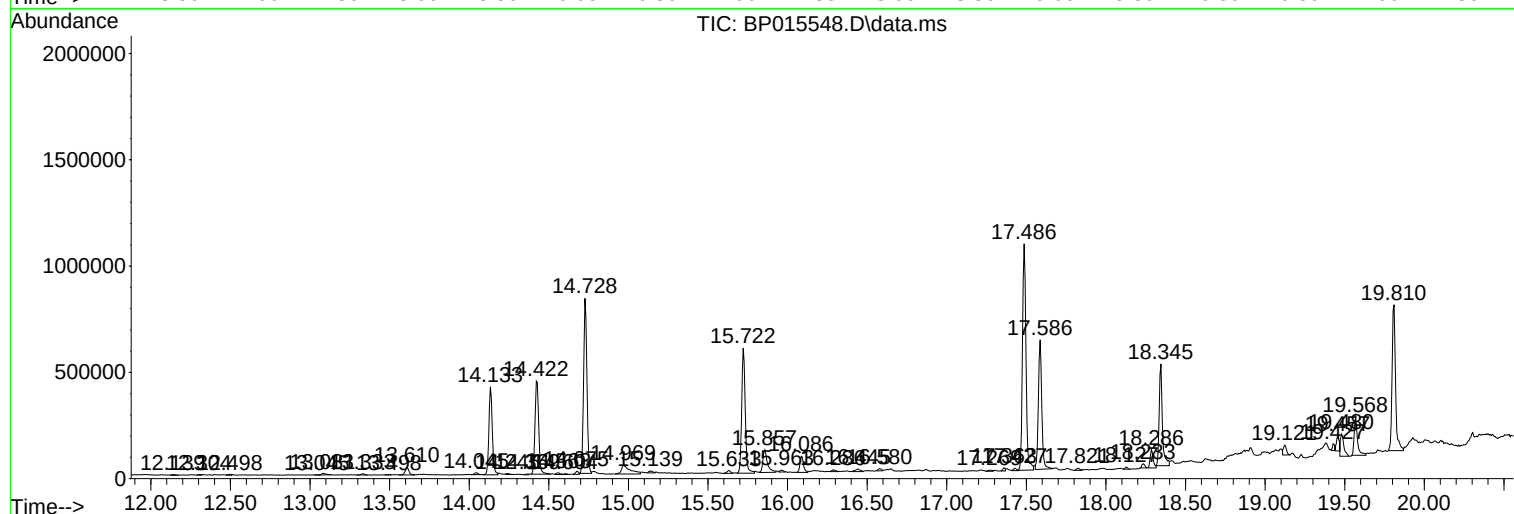
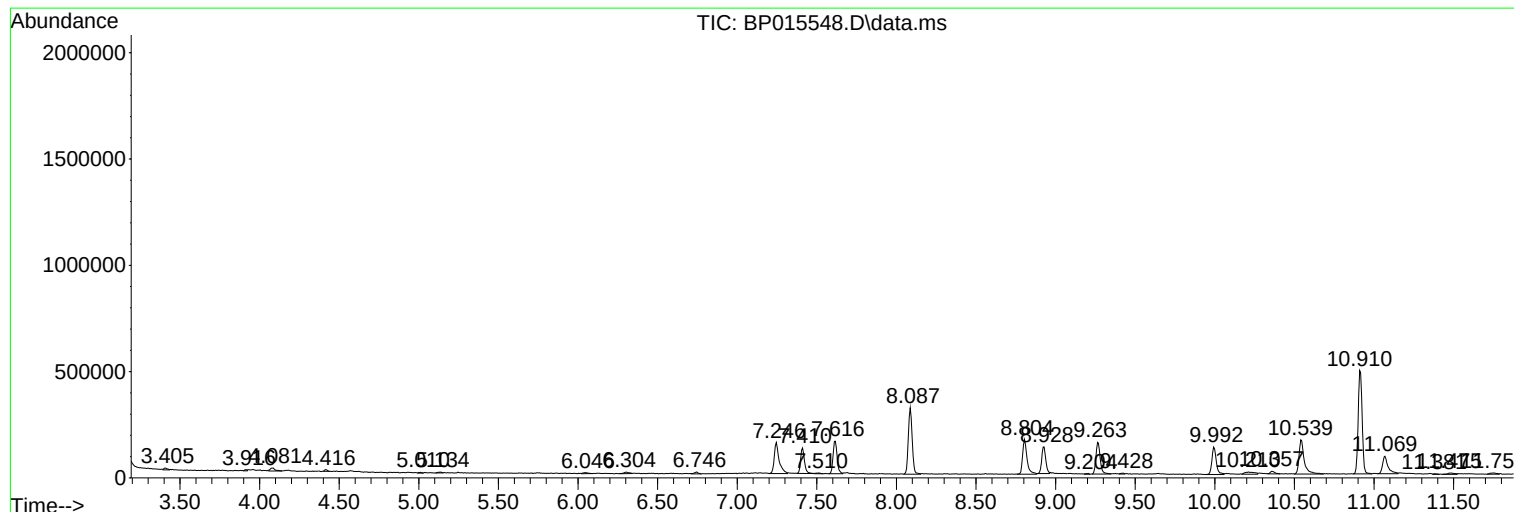
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
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ALS Vial : 43 Sample Multiplier: 1

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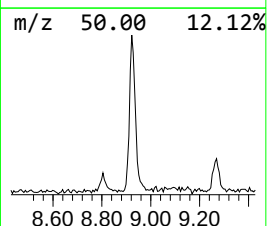
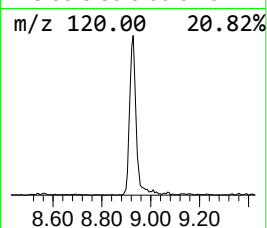
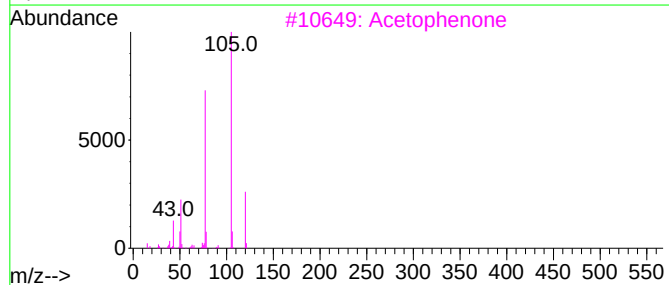
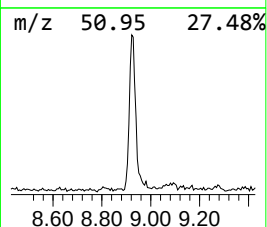
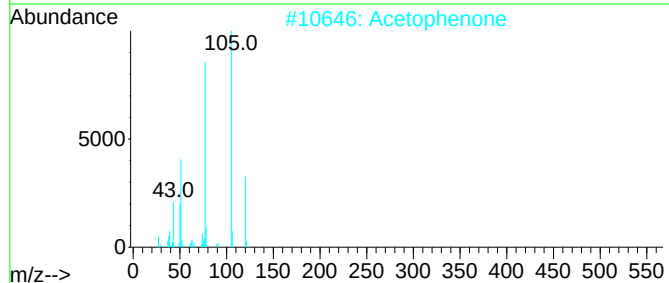
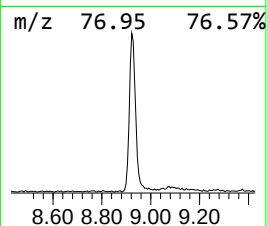
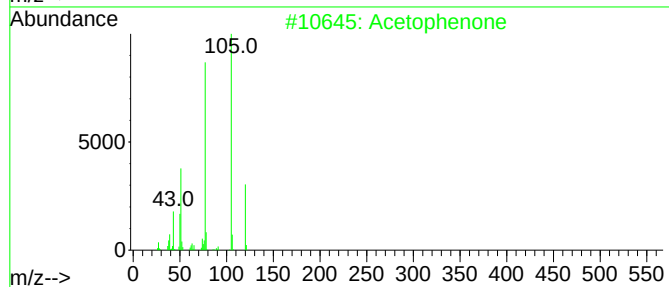
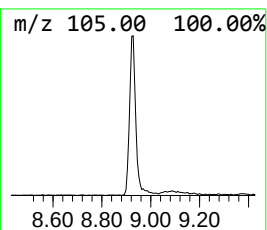
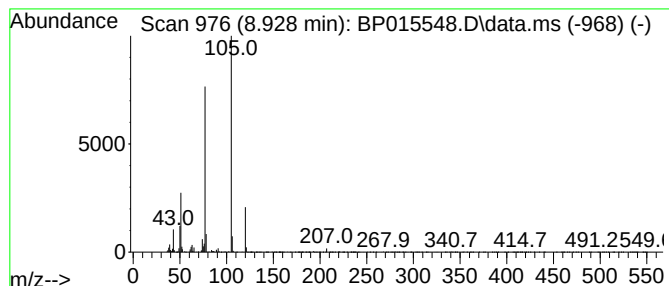
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.928	8.59 ng/ul	223400	1,4-Dichlorobenzene-d4	8.087

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	97
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	90



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Misc :
ALS Vial : 43 Sample Multiplier: 1

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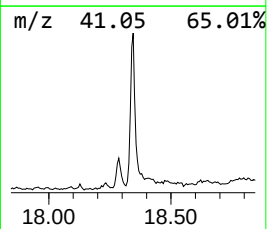
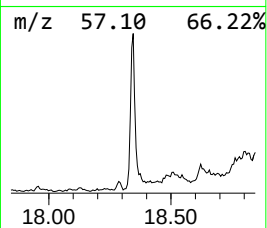
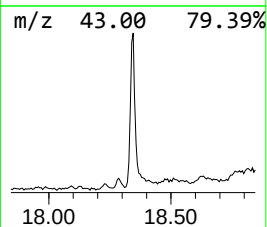
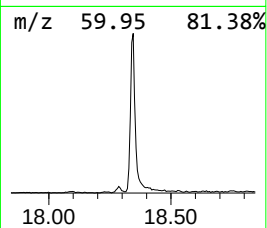
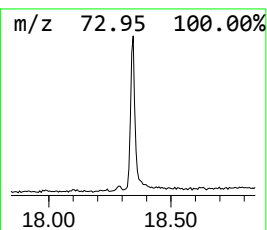
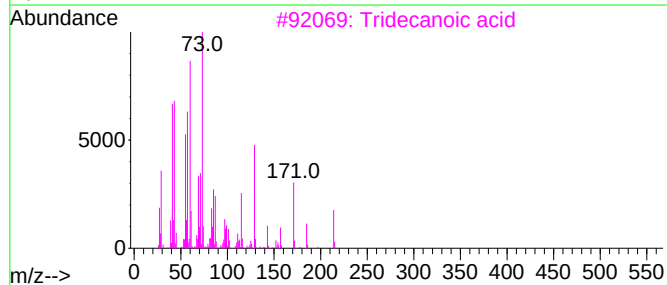
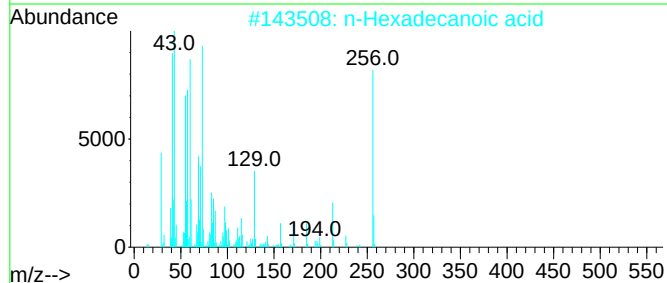
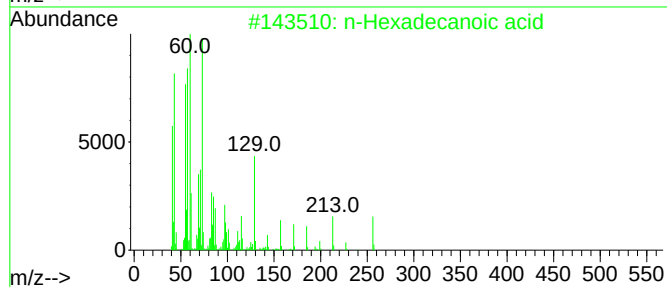
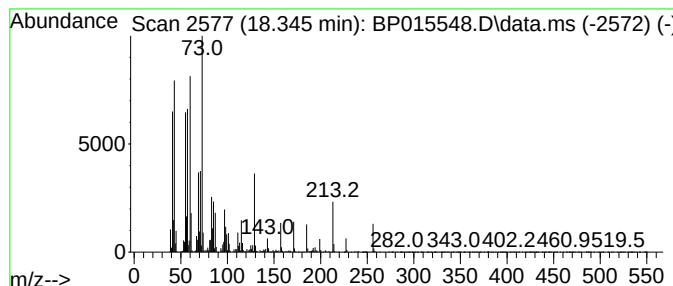
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	8.12 ng/ul	656348	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			Tridecanoic acid	214	C13H26O2	000638-53-9	95
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	94
5			Tridecanoic acid	214	C13H26O2	000638-53-9	93



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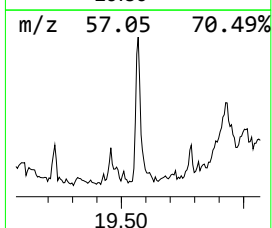
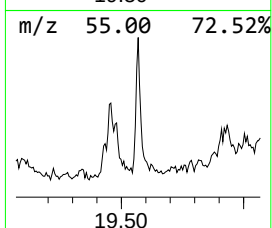
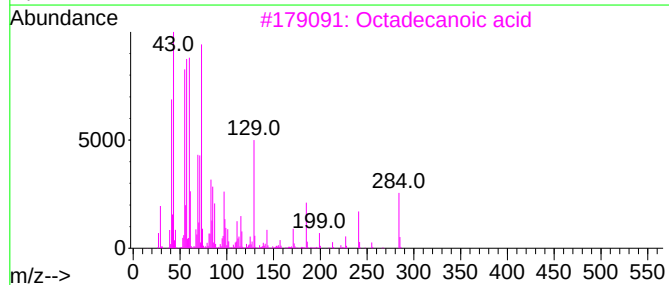
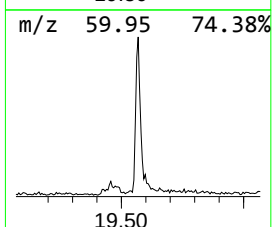
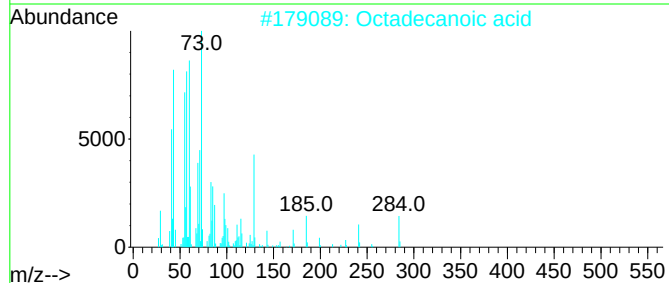
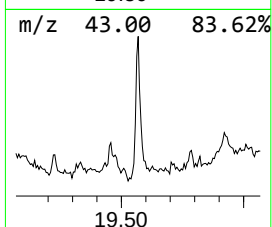
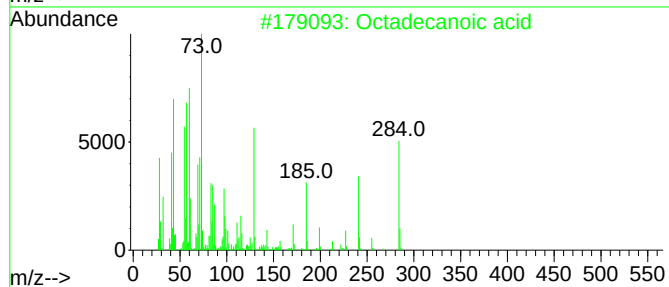
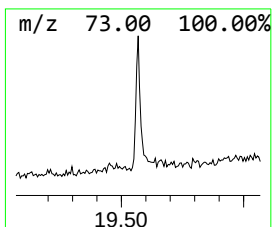
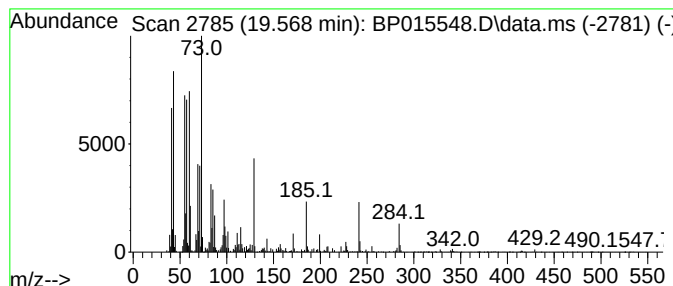
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.568	2.77 ng/ul	253783	Chrysene-d12	21.568

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	99
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	91
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



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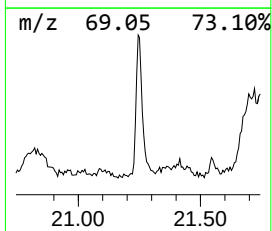
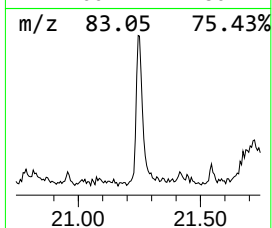
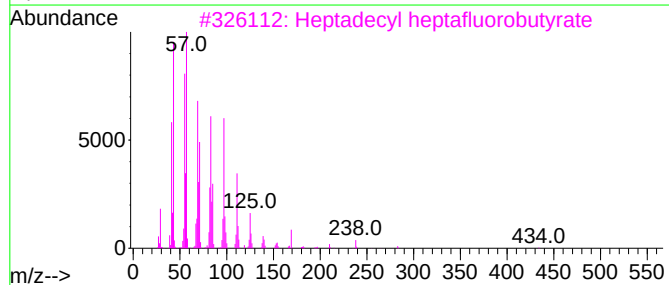
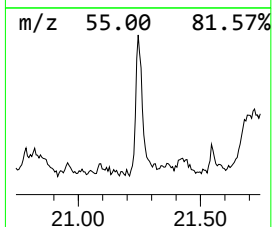
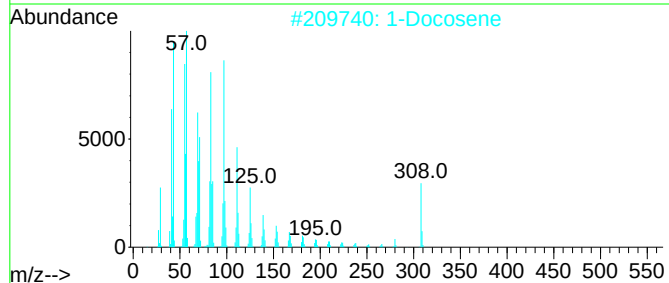
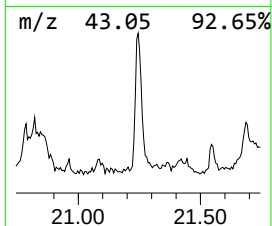
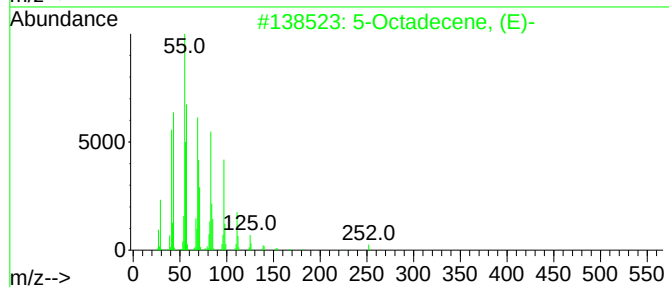
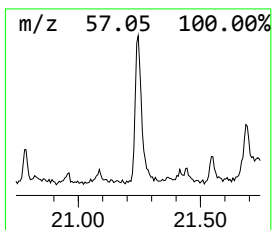
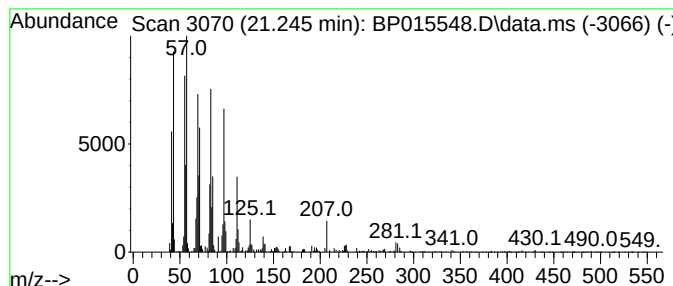
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.245	4.73 ng/ul	433323	Chrysene-d12	21.568

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Octadecene, (E)-	252	C18H36	007206-21-5	95
2			1-Docosene	308	C22H44	001599-67-3	93
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	93
4			1-Hexacosene	364	C26H52	018835-33-1	93
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
Operator : MA/JU
Sample : 03088-07
Misc :
ALS Vial : 43 Sample Multiplier: 1

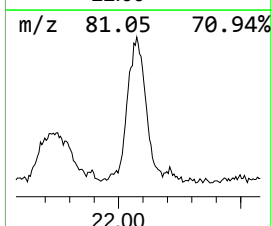
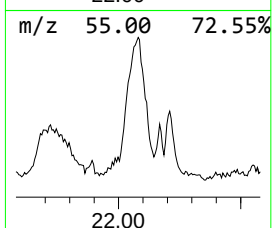
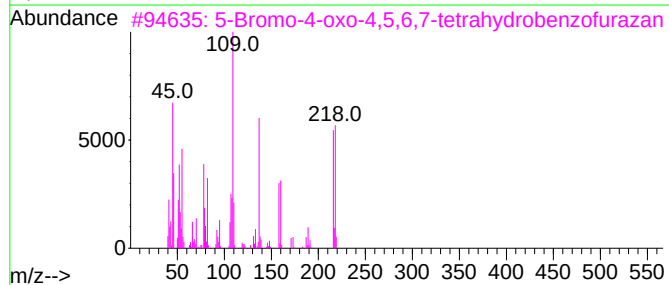
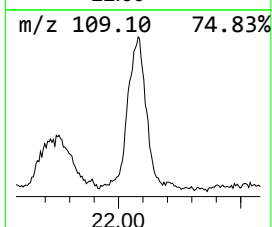
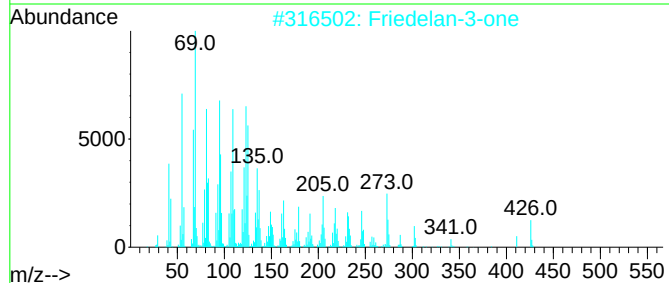
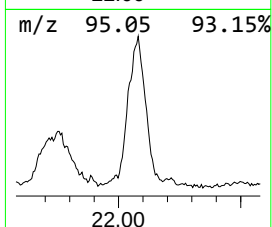
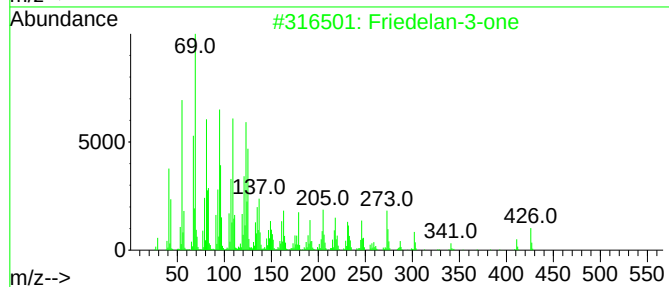
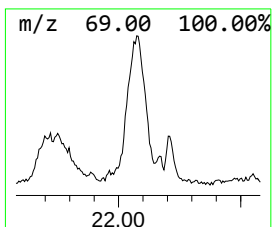
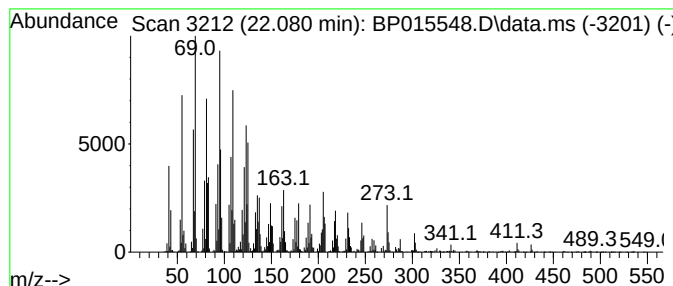
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.080	24.60 ng/ul	2252440	Chrysene-d12		21.568	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Friedelan-3-one		426	C30H50O	000559-74-0	99
2	Friedelan-3-one		426	C30H50O	000559-74-0	95
3	5-Bromo-4-oxo-4,5,6,7-tetrahydro...		216	C6H5BrN2O2	300574-36-1	60
4	Costunolide		232	C15H20O2	000553-21-9	50
5	3-Cyclopentylpropionic acid, but...		194	C12H18O2	1000292-46-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
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Sample : 03088-07
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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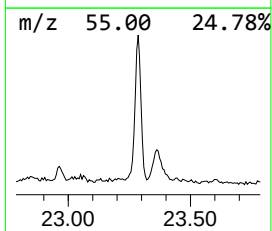
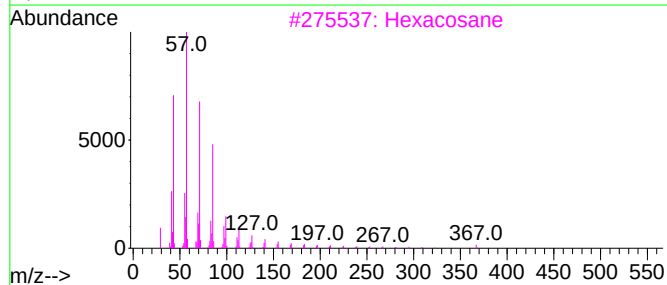
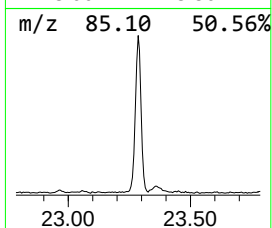
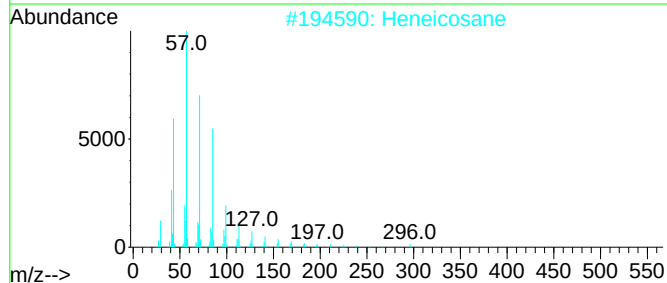
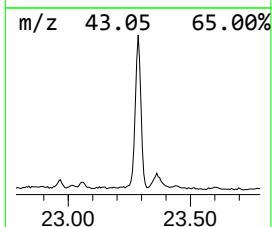
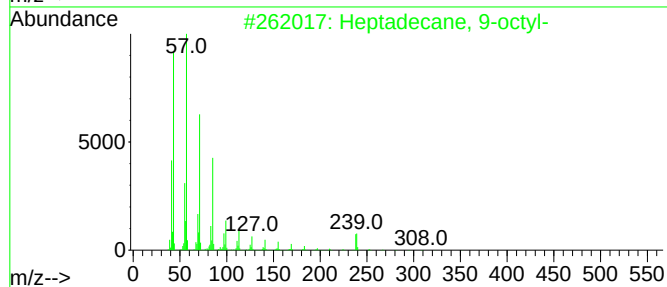
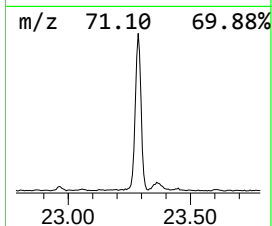
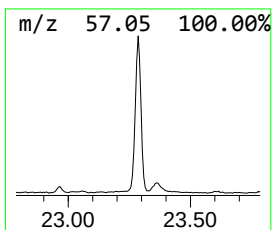
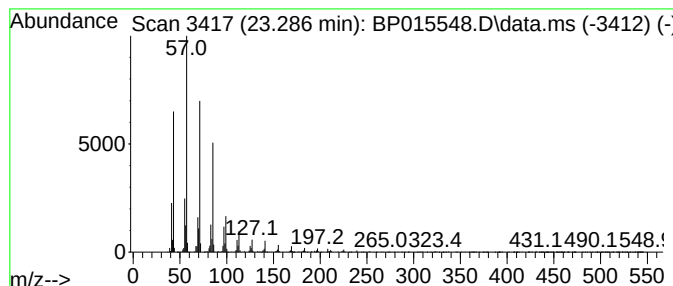
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.286	15.14 ng/ul	1242370	Perylene-d12	24.115

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	96
2			Heneicosane	296	C21H44	000629-94-7	91
3			Hexacosane	366	C26H54	000630-01-3	91
4			Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
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Sample : 03088-07
Misc :
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Instrument :
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ClientSampleId :
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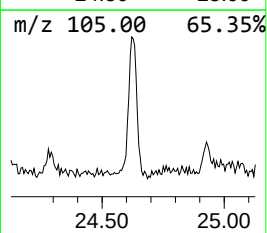
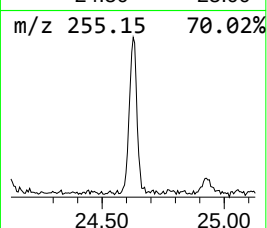
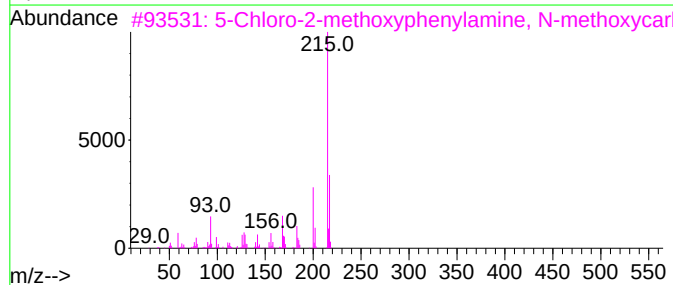
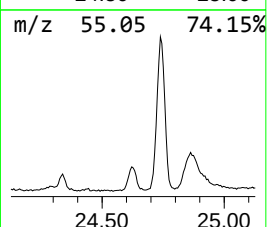
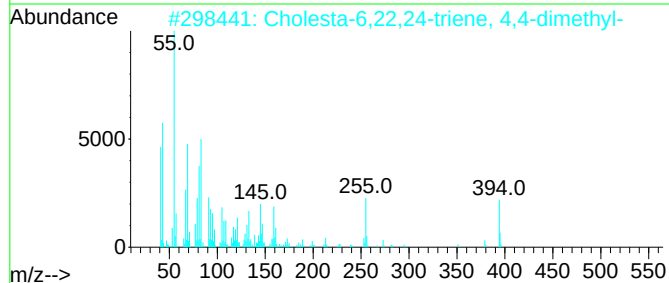
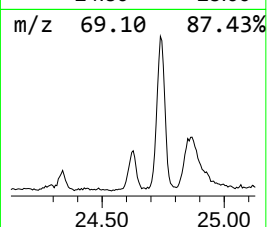
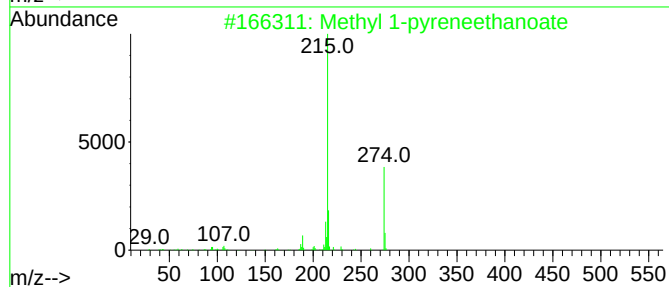
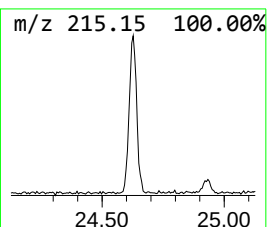
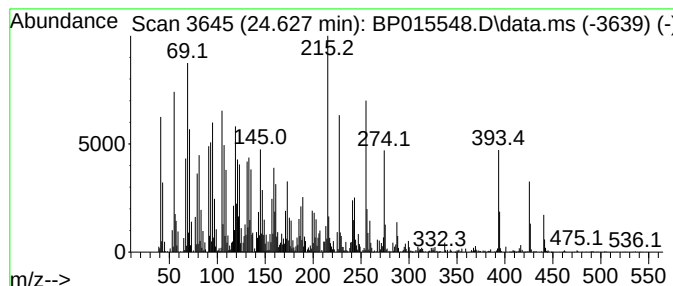
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.627	11.29 ng/ul	926597	Perylene-d12	24.115

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	35
2			Cholesta-6,22,24-triene, 4,4-dim...	394	C29H46	1000128-66-9	35
3			5-Chloro-2-methoxyphenylamine, N...	215	C9H10ClNO3	1000343-98-2	25
4			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	25
5			Stigmasta-5,22-diene, 3-methoxy-...	426	C30H50O	010453-25-5	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
Operator : MA/JU
Sample : 03088-07
Misc :
ALS Vial : 43 Sample Multiplier: 1

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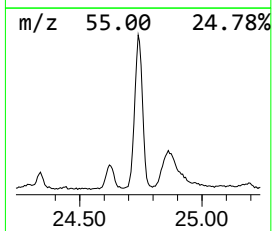
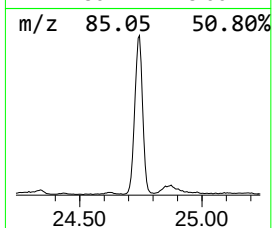
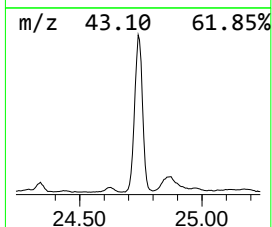
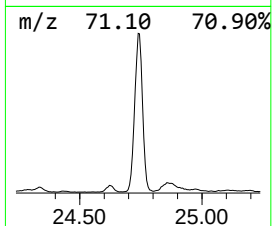
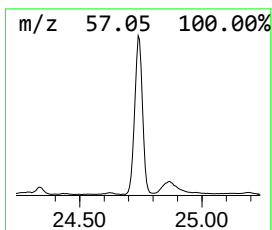
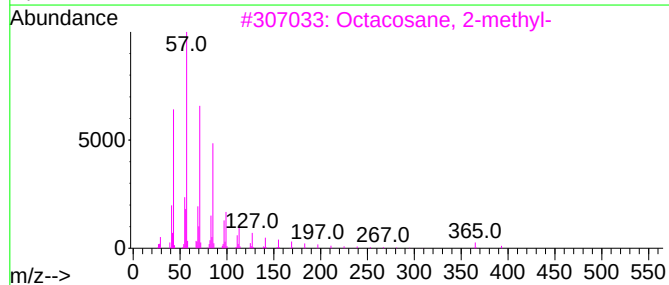
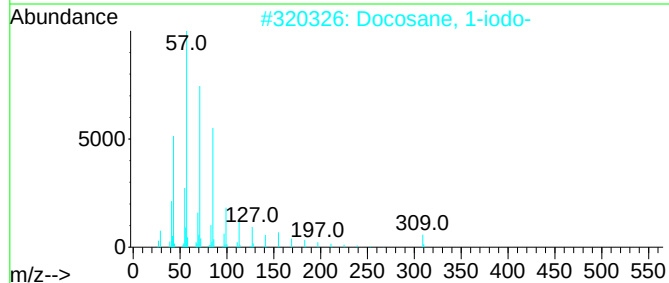
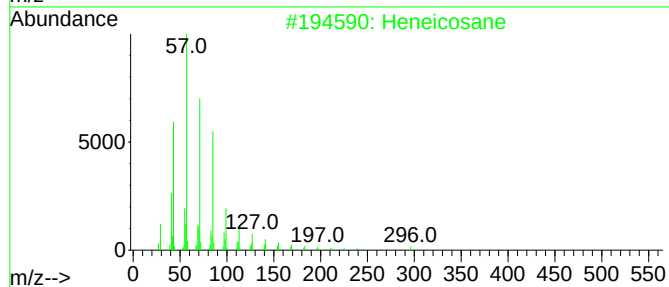
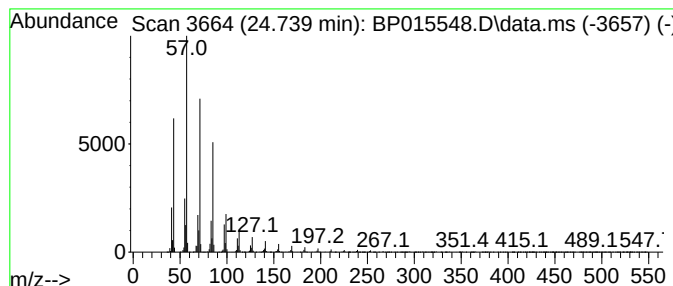
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.739	44.75 ng/ul	3672750	Perylene-d12	24.115

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Octacosane	394	C28H58	000630-02-4	91
5			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
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Sample : 03088-07
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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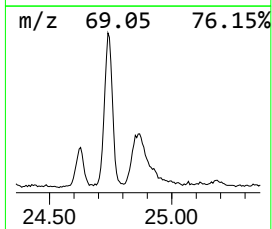
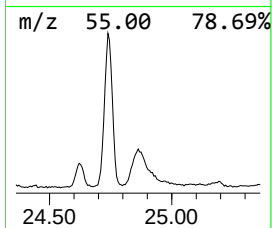
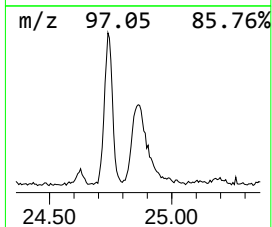
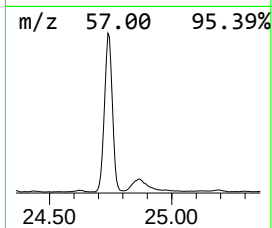
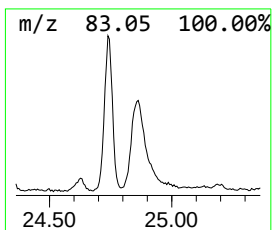
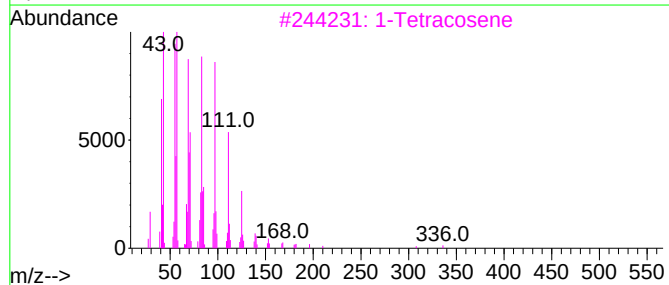
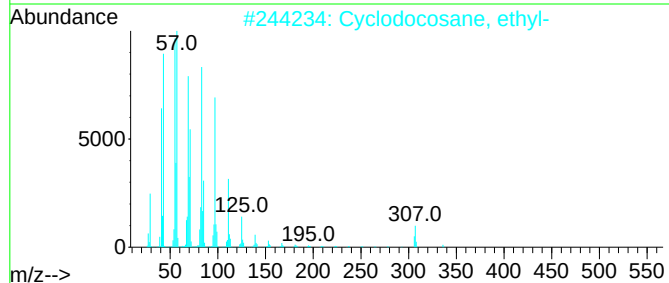
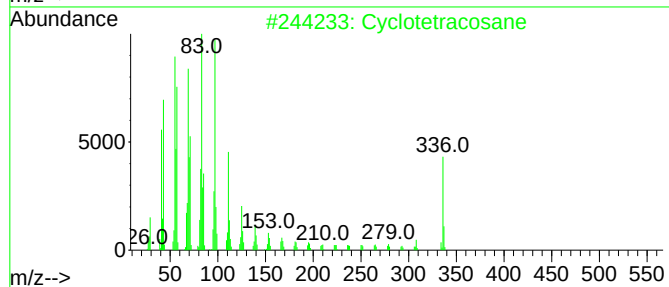
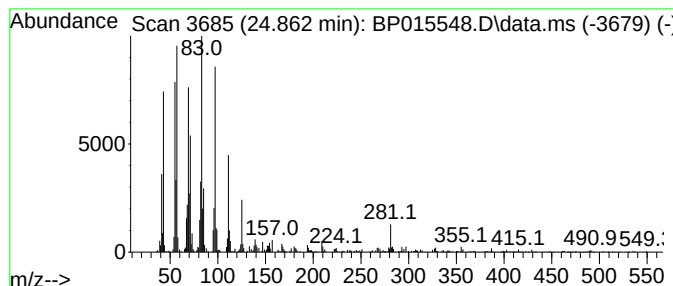
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Cyclic24.862 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.862	6.86 ng/ul	562755	Perylene-d12	24.115

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	96
2			Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	94
3			1-Tetracosene	336	C24H48	010192-32-2	91
4			1-Tricosene	322	C23H46	018835-32-0	90
5			Octacosanol	410	C28H58O	000557-61-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
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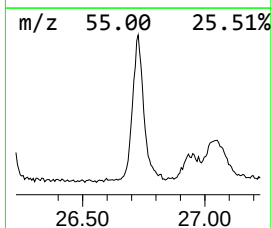
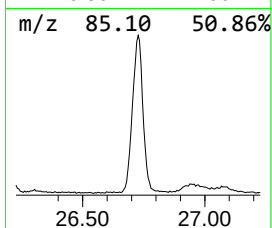
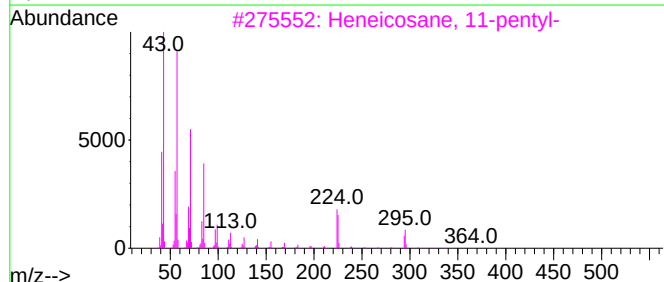
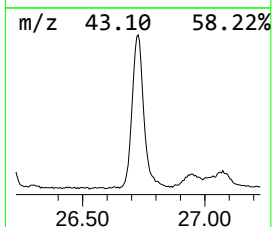
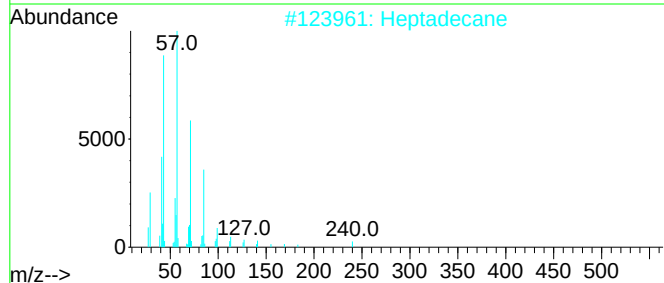
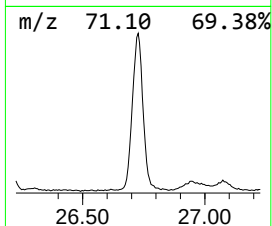
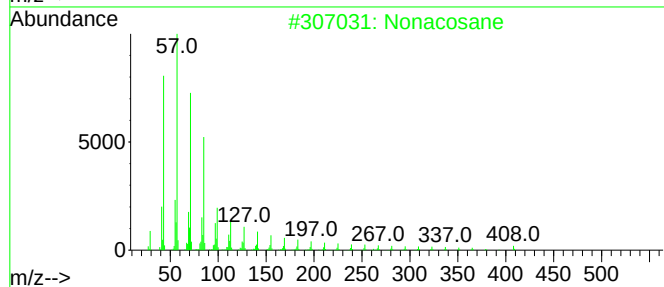
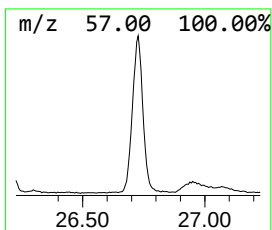
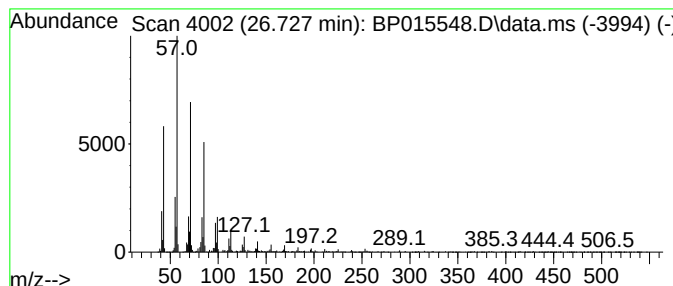
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.727	31.02 ng/ul	2545650	Perylene-d12	24.115

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosane	408	C29H60	000630-03-5	97
2			Heptadecane	240	C17H36	000629-78-7	94
3			Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	94
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	93
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061423\
Data File : BP015548.D
Acq On : 15 Jun 2023 08:53
Operator : MA/JU
Sample : 03088-07
Misc :
ALS Vial : 43 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\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Aceto phenone	8.928	8.6	ng/ul	223400	1	8.087	520381	20.0
n-Hexadecanoic ...	18.345	8.1	ng/ul	656348	4	17.486	1615740	20.0
Octadecanoic acid	19.568	2.8	ng/ul	253783	5	21.568	1831080	20.0
(DEL) Alkane: S...	21.245	4.7	ng/ul	433323	5	21.568	1831080	20.0
5-Bromo-4-oxo-4...	22.080	24.6	ng/ul	2252440	5	21.568	1831080	20.0
(DEL) Alkane: S...	23.286	15.1	ng/ul	1242370	6	24.115	1641560	20.0
unknown-01	24.627	11.3	ng/ul	926597	6	24.115	1641560	20.0
(DEL) Alkane: S...	24.739	44.8	ng/ul	3672750	6	24.115	1641560	20.0
(DEL) Alkane: C...	24.862	6.9	ng/ul	562755	6	24.115	1641560	20.0
(DEL) Alkane: S...	26.727	31.0	ng/ul	2545650	6	24.115	1641560	20.0