

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
 Data File : BP015817.D
 Acq On : 29 Jun 2023 19:12
 Operator : MA/JU
 Sample : 03143-21
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFV8

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: BP015817.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.381	30	33	40	rVB2	35459	54348	1.41%	0.096%
2	3.887	117	119	123	rBV3	14792	25798	0.67%	0.046%
3	3.928	123	126	132	rVV	47843	69034	1.79%	0.122%
4	3.999	136	138	142	rVV5	7957	13459	0.35%	0.024%
5	4.046	142	146	153	rVB	47674	74793	1.94%	0.132%
6	4.152	160	164	169	rVB	27835	38900	1.01%	0.069%
7	4.387	201	204	209	rVB3	14559	18861	0.49%	0.033%
8	6.017	477	481	487	rBV3	22223	38099	0.99%	0.067%
9	6.699	591	597	604	rVB3	25429	49200	1.28%	0.087%
10	6.911	629	633	639	rBV8	8695	13886	0.36%	0.025%
11	7.240	682	689	708	rBV	301957	884683	22.93%	1.565%
12	7.387	708	714	733	rVB	346922	696547	18.06%	1.232%
13	7.593	742	749	756	rBV	450583	858665	22.26%	1.519%
14	7.646	756	758	764	rVB	48388	63419	1.64%	0.112%
15	8.005	815	819	822	rBV5	11685	18714	0.49%	0.033%
16	8.058	822	828	843	rVV	1053107	1987244	51.51%	3.516%
17	8.316	867	872	878	rVB9	7676	14341	0.37%	0.025%
18	8.705	934	938	945	rBV9	7300	13927	0.36%	0.025%
19	8.799	946	954	969	rBV	355051	964875	25.01%	1.707%
20	8.916	969	974	992	rVB	145230	340894	8.84%	0.603%
21	9.252	1024	1031	1047	rBV	372649	877994	22.76%	1.553%
22	9.593	1086	1089	1094	rVB3	9585	15413	0.40%	0.027%
23	9.981	1147	1155	1176	rBV	241063	715416	18.55%	1.266%
24	10.275	1196	1205	1208	rBV8	17690	46449	1.20%	0.082%
25	10.381	1219	1223	1228	rVB6	11529	20376	0.53%	0.036%
26	10.558	1245	1253	1272	rBV	283791	980985	25.43%	1.736%
27	10.840	1295	1301	1302	rBV2	23928	35067	0.91%	0.062%
28	10.887	1302	1309	1328	rVV	1475591	2956637	76.64%	5.231%
29	11.016	1328	1331	1336	rVV4	14571	20648	0.54%	0.037%
30	11.081	1336	1342	1360	rVV3	101826	340654	8.83%	0.603%
31	11.787	1456	1462	1469	rVB3	8728	20428	0.53%	0.036%
32	11.881	1473	1478	1483	rBV2	28749	48549	1.26%	0.086%
33	12.281	1541	1546	1555	rBV2	30859	62168	1.61%	0.110%
34	12.416	1564	1569	1571	rBV4	8670	15099	0.39%	0.027%
35	12.499	1578	1583	1589	rVB8	12473	26182	0.68%	0.046%
36	12.569	1589	1595	1604	rBV	21314	50070	1.30%	0.089%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	12.781	1626	1631	1636	rBV3	20907	39416	1.02%	0.070%
38	13.075	1679	1681	1685	rVV5	10342	15661	0.41%	0.028%
39	13.216	1700	1705	1712	rBV3	24692	39911	1.03%	0.071%
40	13.504	1749	1754	1760	rBV	42348	68997	1.79%	0.122%

41	13.587	1760	1768	1775	rVV2	49143	105441	2.73%	0.187%
42	13.681	1778	1784	1789	rVB9	9123	18433	0.48%	0.033%
43	13.899	1812	1821	1826	rBV5	15751	42948	1.11%	0.076%
44	13.975	1828	1834	1837	rVV7	7210	13835	0.36%	0.024%
45	14.028	1837	1843	1848	rVV3	28679	59659	1.55%	0.106%

46	14.122	1848	1859	1880	rVV	867641	1689152	43.79%	2.989%
47	14.269	1880	1884	1893	rVB4	14750	38649	1.00%	0.068%
48	14.404	1900	1907	1921	rBV	1168914	1959790	50.80%	3.467%
49	14.510	1923	1925	1931	rVB7	15833	23099	0.60%	0.041%
50	14.575	1931	1936	1942	rVB	48985	68298	1.77%	0.121%

51	14.651	1944	1949	1952	rBV6	12288	18566	0.48%	0.033%
52	14.704	1952	1958	1974	rBV	2296953	3679607	95.38%	6.510%
53	15.016	2005	2011	2018	rBV5	63526	198432	5.14%	0.351%
54	15.122	2026	2029	2034	rVB4	28331	44654	1.16%	0.079%
55	15.187	2037	2040	2047	rVB7	22090	45109	1.17%	0.080%

56	15.328	2059	2064	2068	rBV7	12024	24155	0.63%	0.043%
57	15.381	2070	2073	2085	rVB9	12265	27331	0.71%	0.048%
58	15.475	2085	2089	2094	rBV7	20764	42299	1.10%	0.075%
59	15.528	2094	2098	2102	rVB	51902	70669	1.83%	0.125%
60	15.704	2121	2128	2142	rBV	1248999	2294279	59.47%	4.059%

61	15.881	2152	2158	2164	rBV	148250	337742	8.76%	0.598%
62	16.075	2185	2191	2201	rVB2	151173	265422	6.88%	0.470%
63	16.216	2211	2215	2217	rBV4	10243	14899	0.39%	0.026%
64	16.310	2228	2231	2238	rVB7	15201	31471	0.82%	0.056%
65	16.387	2240	2244	2246	rBV2	62607	85162	2.21%	0.151%

66	16.557	2269	2273	2279	rBV7	25248	44378	1.15%	0.079%
67	16.845	2319	2322	2327	rBV6	17340	24301	0.63%	0.043%
68	16.998	2343	2348	2351	rBV7	14508	24477	0.63%	0.043%
69	17.187	2376	2380	2385	rBV3	61609	72304	1.87%	0.128%
70	17.339	2402	2406	2413	rBV4	56816	87848	2.28%	0.155%

71	17.404	2413	2417	2421	rBV3	35696	54252	1.41%	0.096%
72	17.469	2421	2428	2433	rBV	2535320	3857666	100.00%	6.825%
73	17.569	2440	2445	2457	rVB	1141247	1979996	51.33%	3.503%
74	17.792	2481	2483	2489	rVB7	15701	21685	0.56%	0.038%
75	17.922	2498	2505	2510	rBV2	72228	120755	3.13%	0.214%

76	18.098	2533	2535	2539	rVB2	29206	32278	0.84%	0.057%
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77	18.210	2550	2554	2559	rBV6	34238	51053	1.32%	0.090%
78	18.263	2559	2563	2568	rBV	282806	380039	9.85%	0.672%
79	18.322	2568	2573	2581	rBV	1577113	2143661	55.57%	3.793%
80	18.516	2602	2606	2610	rBV3	51188	94197	2.44%	0.167%
81	18.586	2616	2618	2626	rVB	72167	105412	2.73%	0.187%
82	19.198	2718	2722	2728	rBV3	114481	170973	4.43%	0.303%
83	19.304	2736	2740	2744	rBV3	83911	124957	3.24%	0.221%
84	19.404	2754	2757	2759	rBV	108699	121233	3.14%	0.214%
85	19.469	2763	2768	2775	rVB	629842	1028077	26.65%	1.819%
86	19.545	2777	2781	2786	rBV2	398094	550237	14.26%	0.974%
87	19.792	2812	2823	2826	rBV	1651076	2858403	74.10%	5.057%
88	20.269	2901	2904	2908	rBV	241721	303330	7.86%	0.537%
89	20.492	2939	2942	2946	rVB4	130725	142748	3.70%	0.253%
90	20.751	2983	2986	2990	rBV	174828	194349	5.04%	0.344%
91	21.210	3060	3064	3065	rBV2	423358	542600	14.07%	0.960%
92	21.416	3096	3099	3105	rBV	315650	360597	9.35%	0.638%
93	21.551	3117	3122	3127	rBV2	2037289	3241774	84.03%	5.736%
94	22.133	3218	3221	3225	rBV	332294	372309	9.65%	0.659%
95	23.239	3404	3409	3415	rVB	927385	1513024	39.22%	2.677%
96	23.322	3418	3423	3439	rVB2	546004	1572205	40.76%	2.782%
97	24.098	3548	3555	3563	rBV	1325296	3021180	78.32%	5.345%
98	24.563	3630	3634	3644	rVB8	529374	1136645	29.46%	2.011%
99	24.674	3646	3653	3662	rVB	1683625	3699539	95.90%	6.546%
100	26.633	3979	3986	3995	rBV2	904292	2636385	68.34%	4.665%

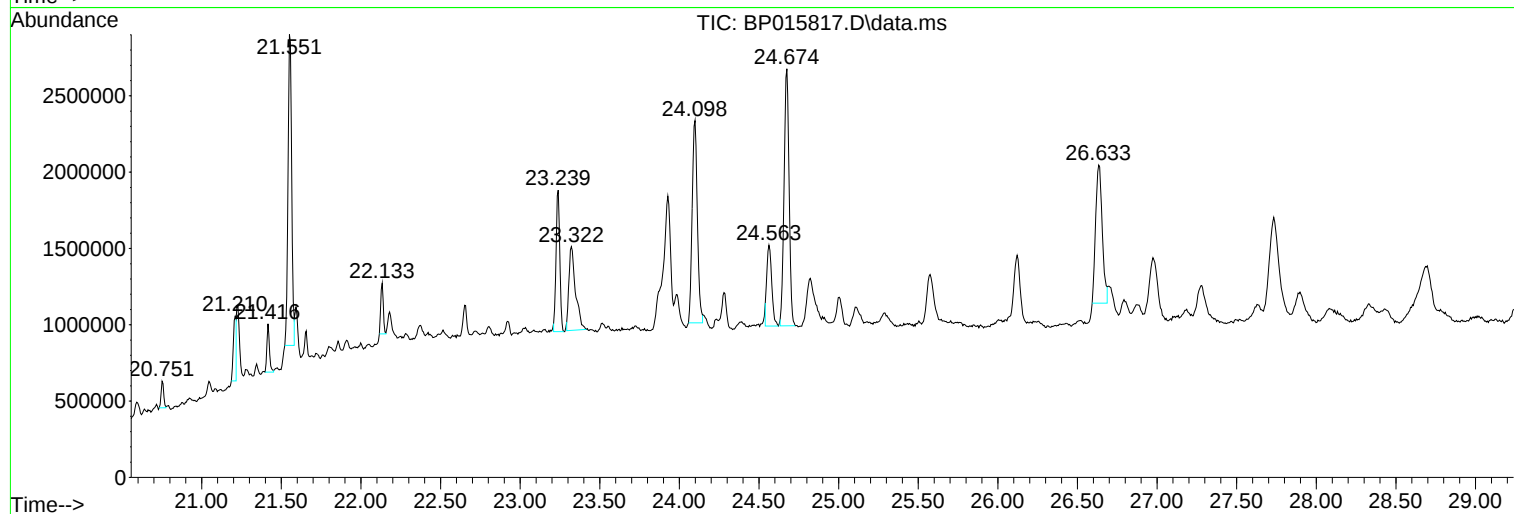
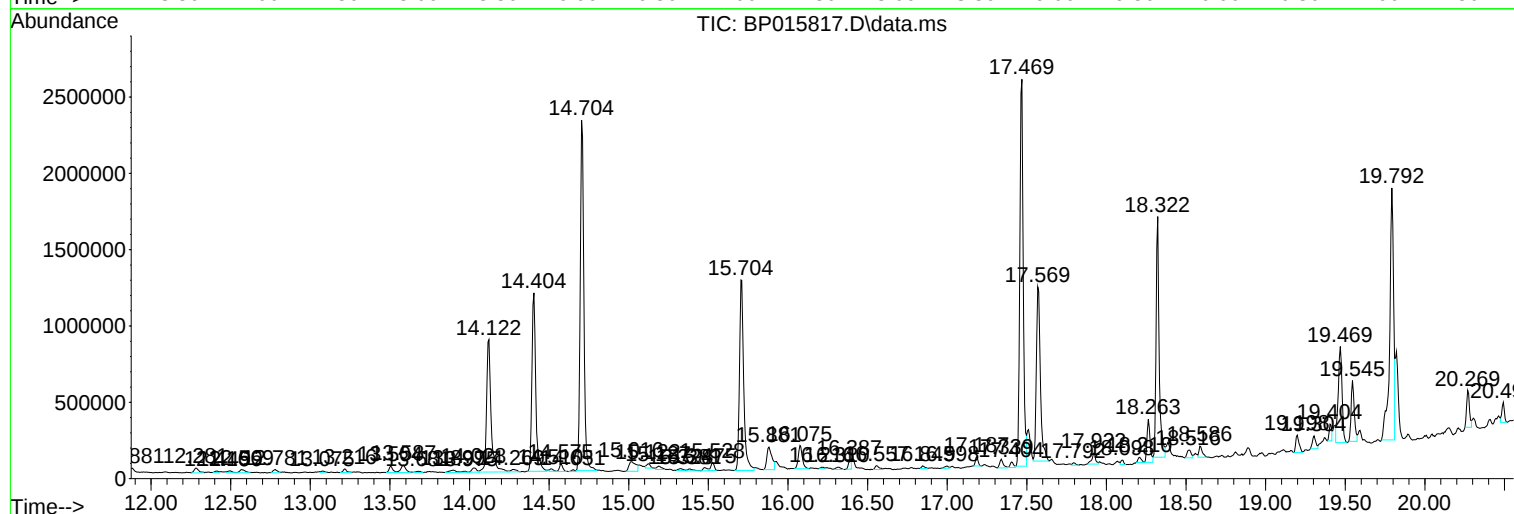
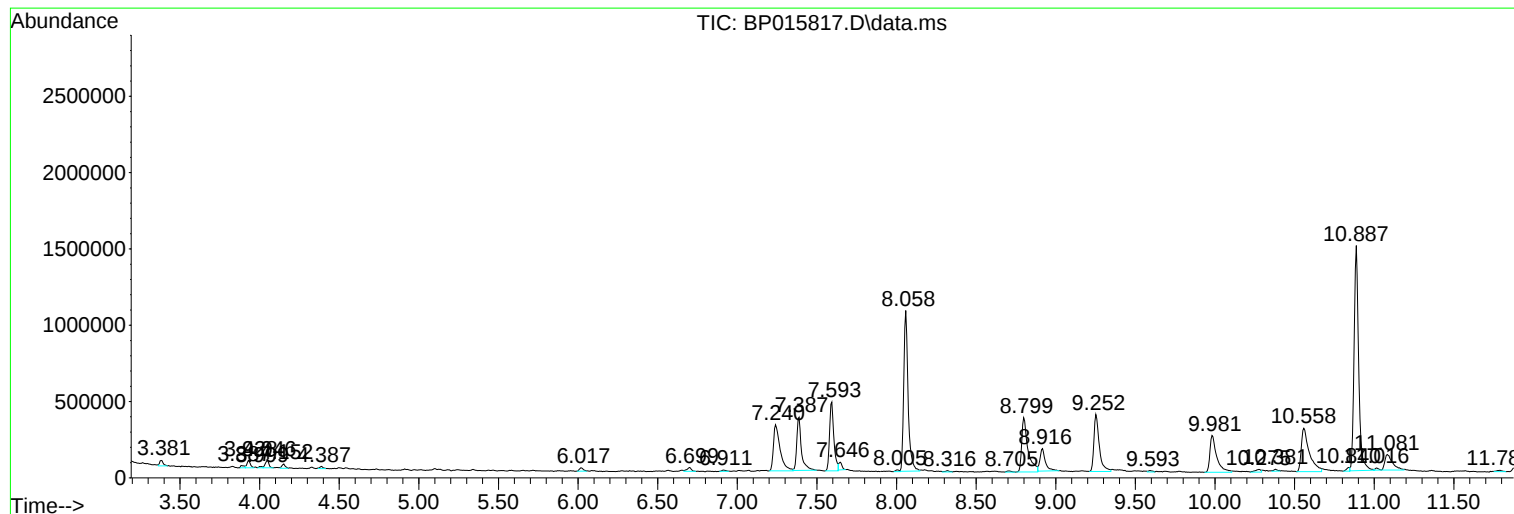
Sum of corrected areas: 56519775

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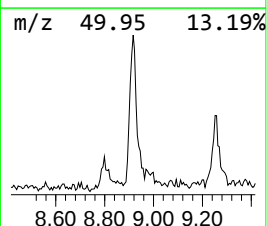
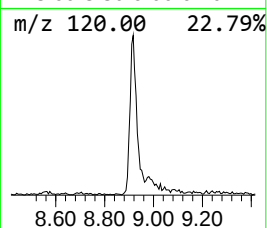
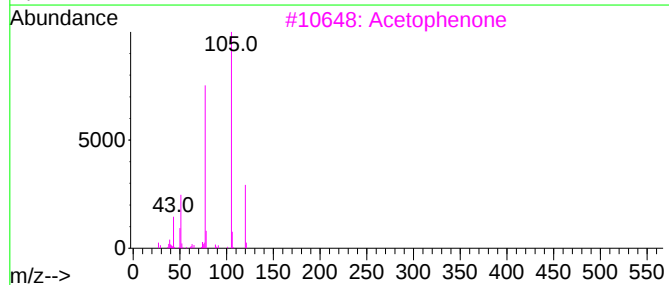
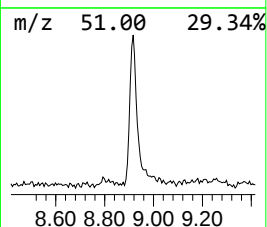
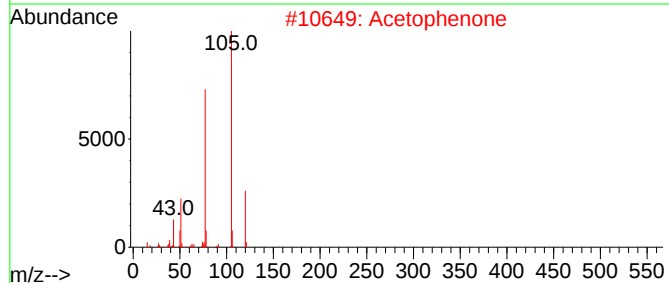
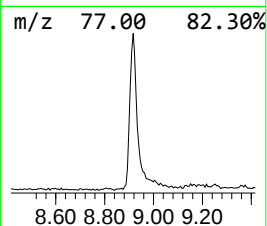
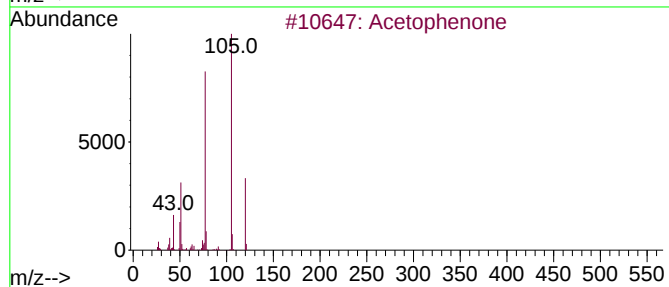
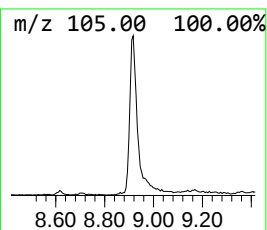
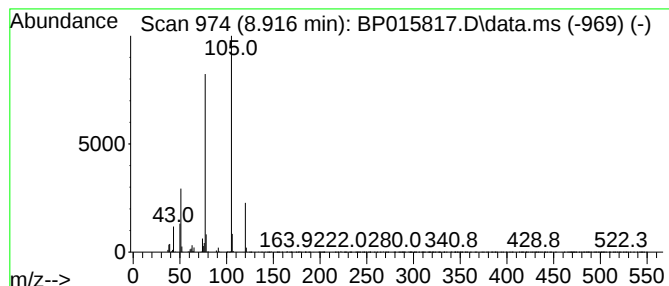
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.916	3.43 ng/ul	340894	1,4-Dichlorobenzene-d4	8.058

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	90
5			1-Butanone, 1-phenyl-	148	C10H12O	000495-40-9	80



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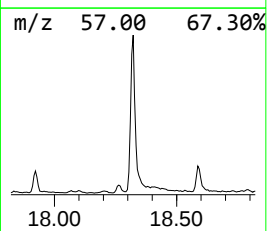
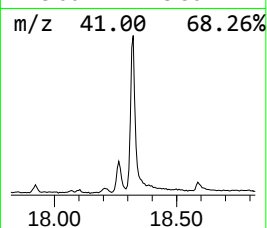
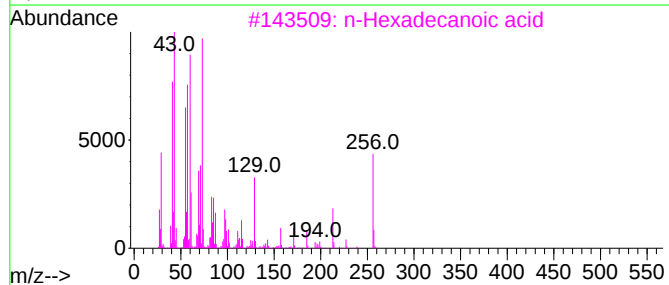
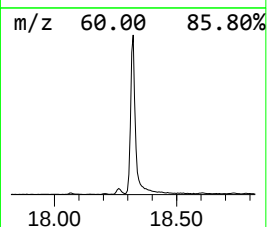
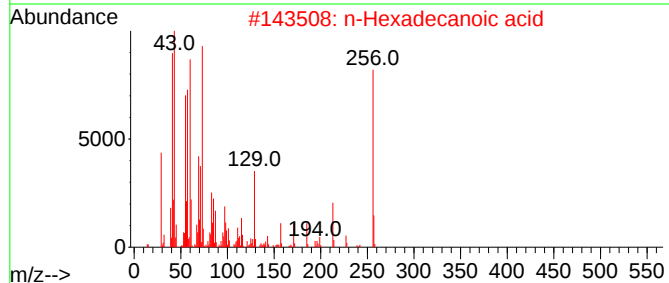
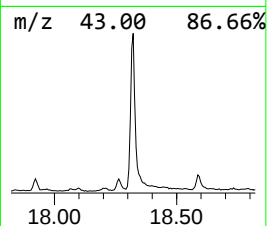
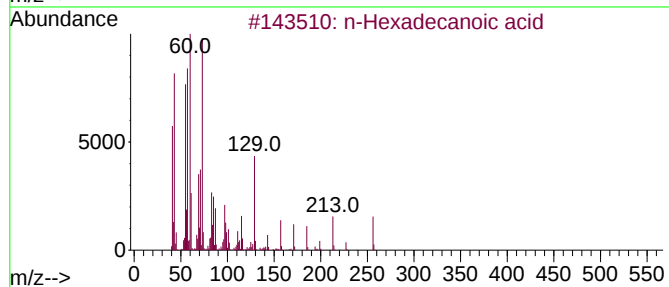
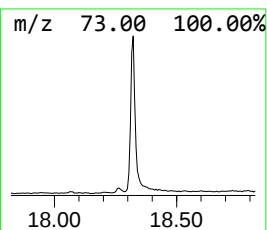
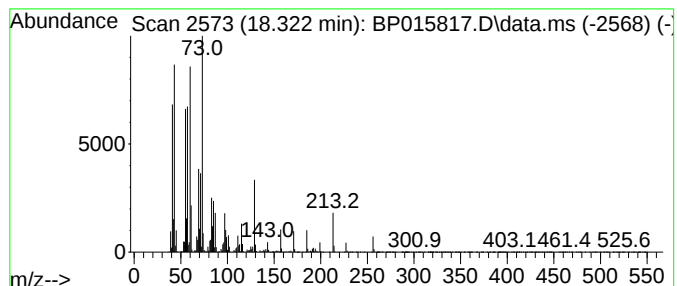
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.322	11.11 ng/ul	2143660	Phenanthrene-d10	17.469

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			Tetradecanoic acid	228	C14H28O2	000544-63-8	95



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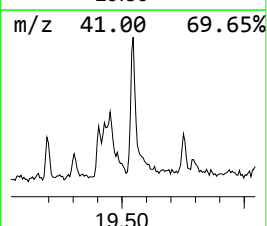
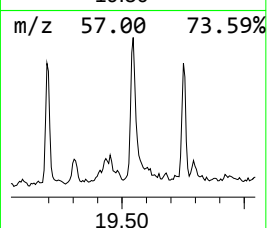
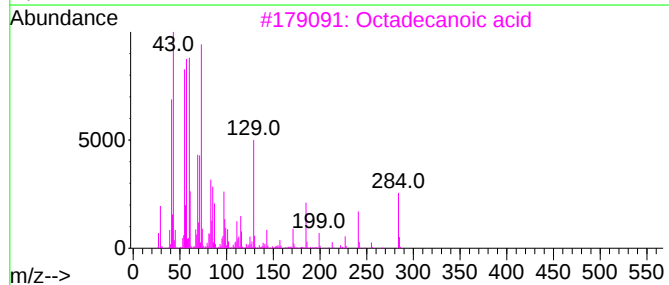
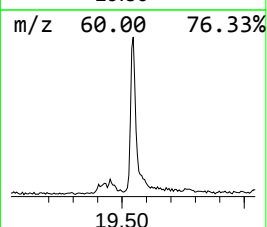
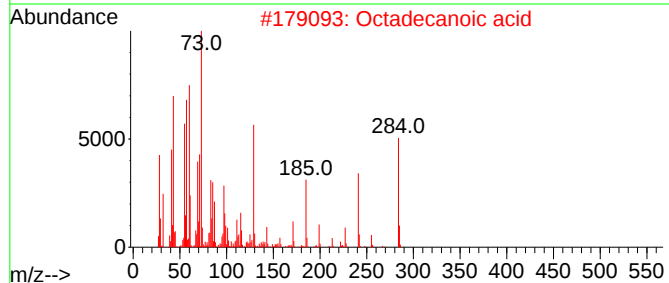
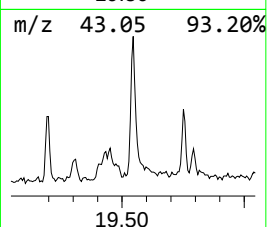
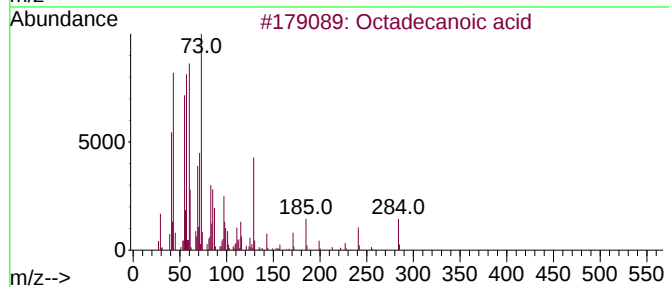
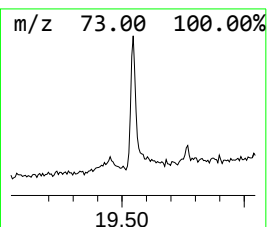
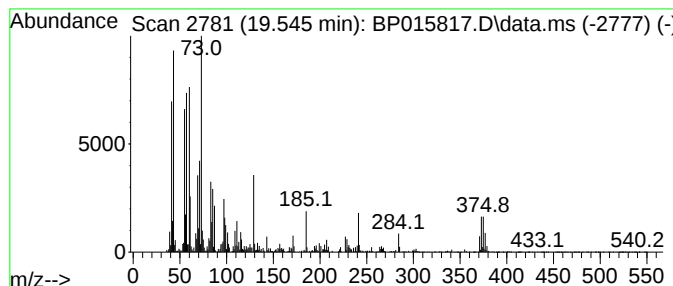
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.545	3.39 ng/ul	550237	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
Operator : MA/JU
Sample : 03143-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFV8

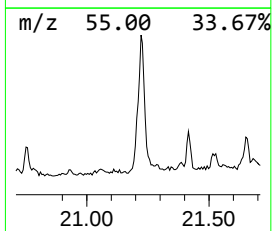
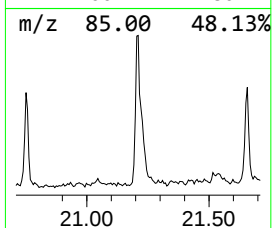
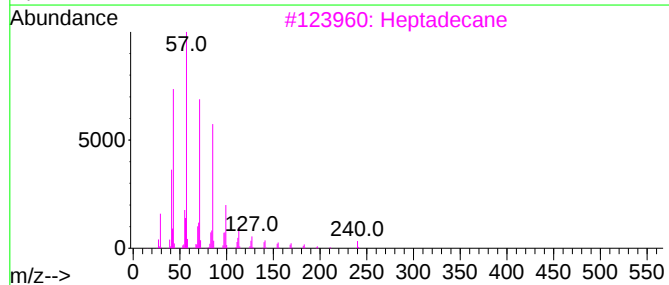
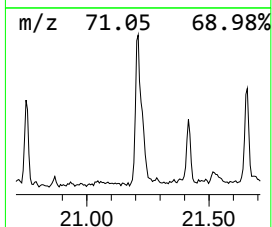
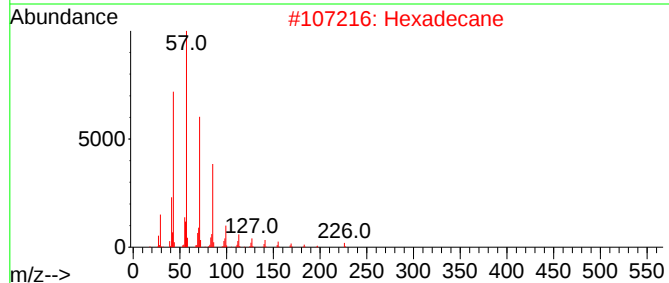
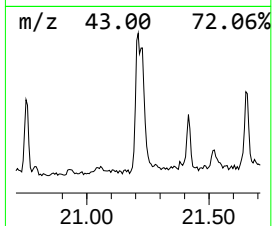
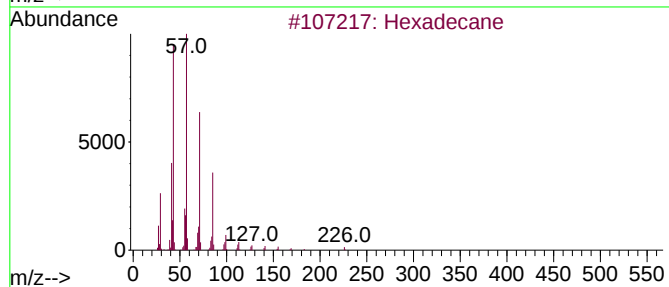
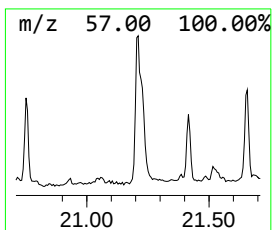
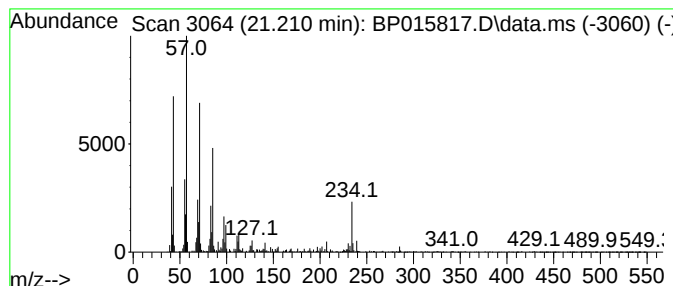
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.210	3.35 ng/ul	542600	Chrysene-d12	21.557

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	93
2		Hexadecane	226	C16H34	000544-76-3	93
3		Heptadecane	240	C17H36	000629-78-7	93
4		Pentacosane	352	C25H52	000629-99-2	92
5		Heptadecane	240	C17H36	000629-78-7	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
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Sample : 03143-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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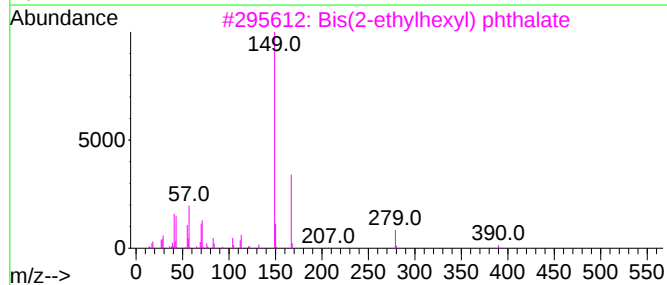
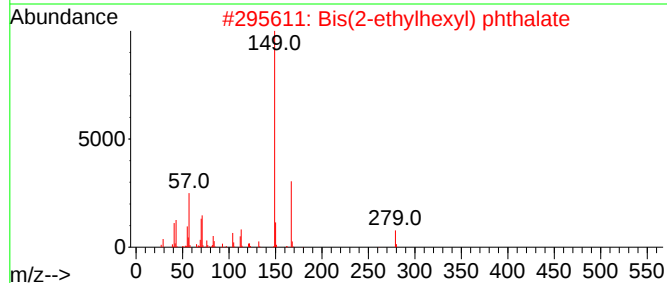
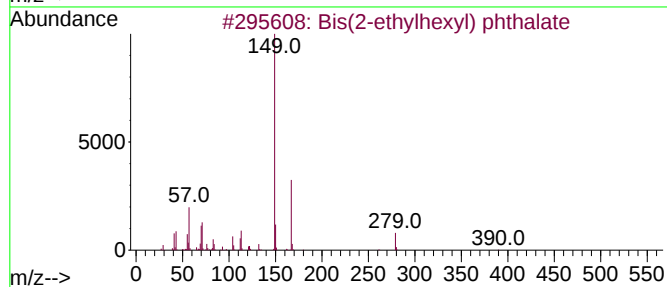
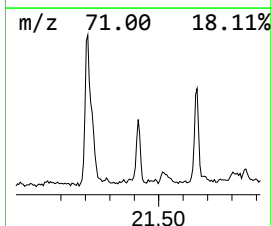
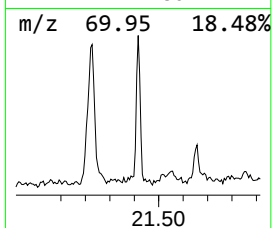
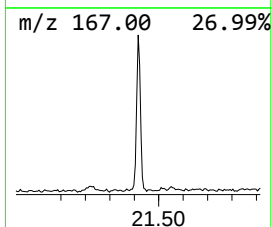
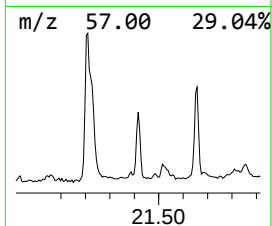
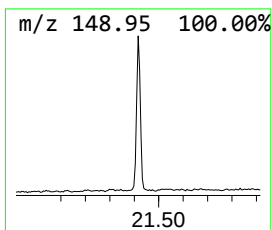
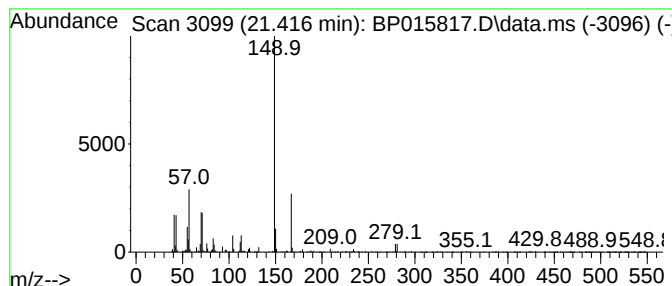
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Bis(2-ethylhexyl) phthalate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.416	2.22 ng/ul	360597	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	93
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	87
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	87
4			Phthalic acid, di(hex-3-yl) ester	334	C20H30O4	1000356-96-7	64
5			Phthalic acid, hept-3-yl isohexy...	348	C21H32O4	1000356-98-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
Operator : MA/JU
Sample : 03143-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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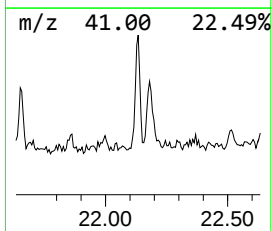
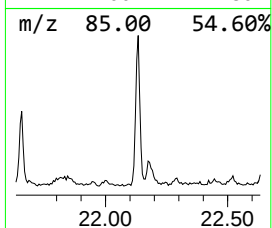
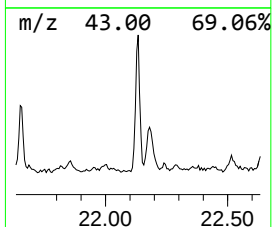
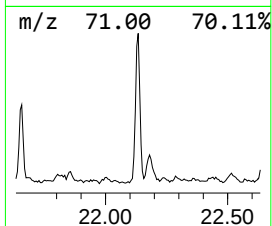
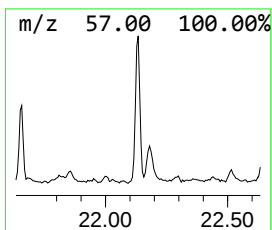
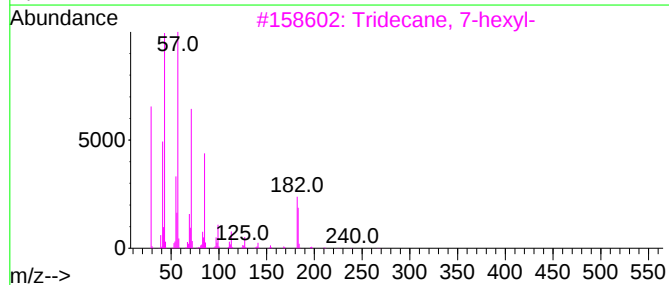
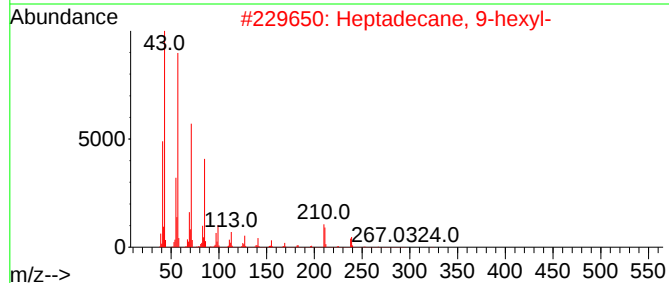
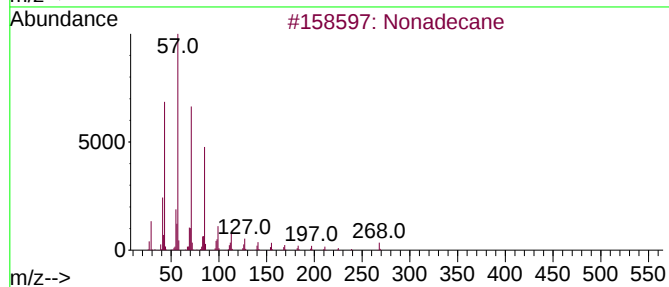
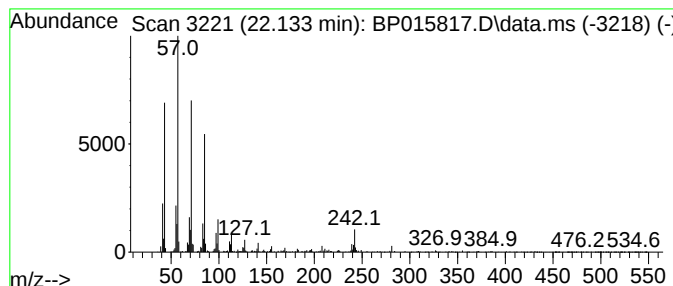
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.133	2.30 ng/ul	372309	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	94
2			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	90
3			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	90
4			Heptadecane	240	C17H36	000629-78-7	83
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
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Sample : 03143-21
Misc :
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Instrument :
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ClientSampleId :
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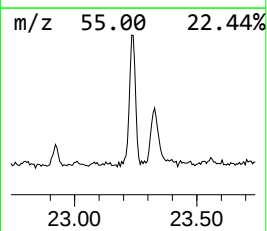
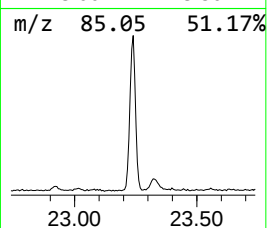
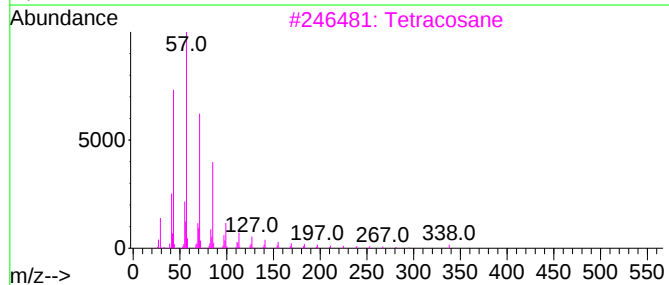
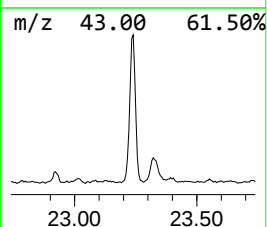
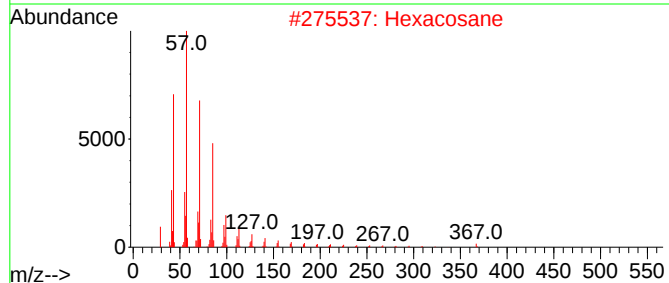
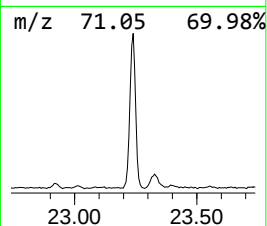
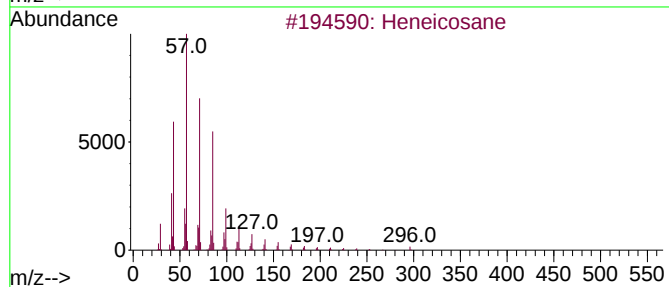
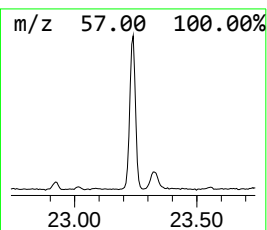
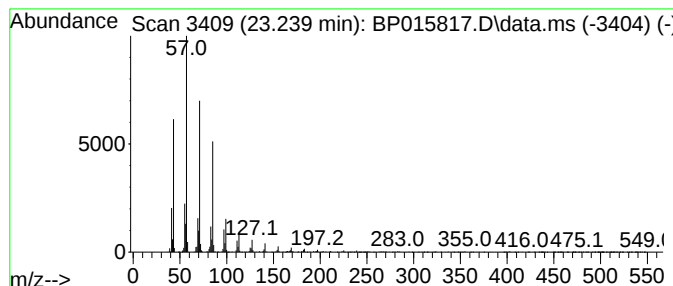
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.239	10.02 ng/ul	1513020	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Hexacosane	366	C26H54	000630-01-3	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5			Tetratetracontane	619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
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Sample : 03143-21
Misc :
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ClientSampleId :
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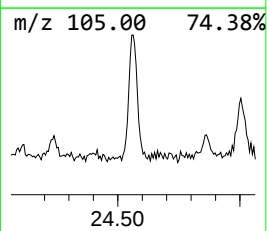
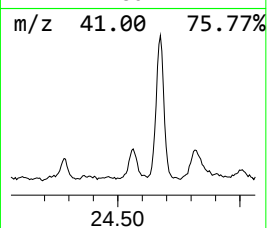
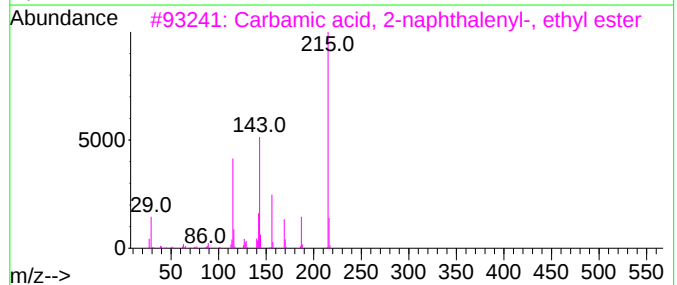
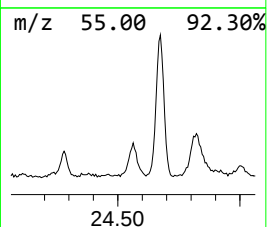
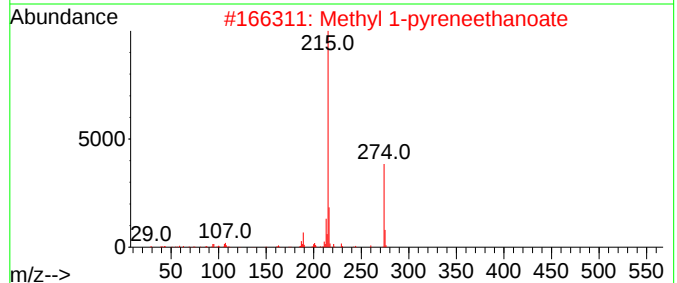
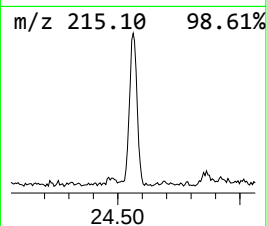
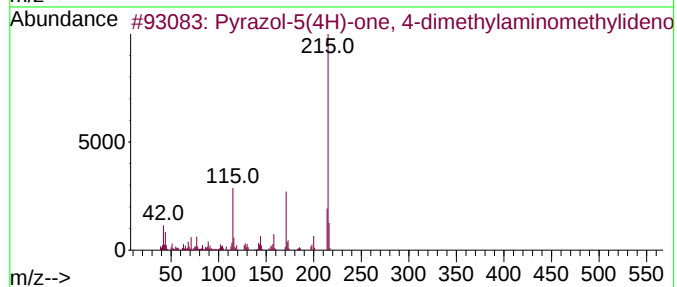
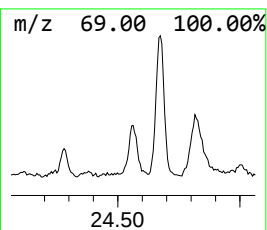
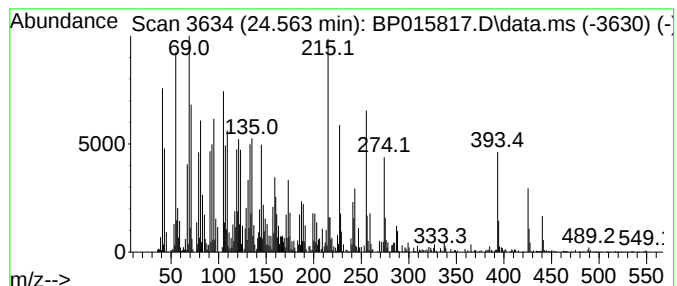
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.563	7.52 ng/ul	1136650	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrazol-5(4H)-one, 4-dimethylami...	215	C12H13N3O	021272-28-6	25
2			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	20
3			Carbamic acid, 2-naphthalenyl-, ...	215	C13H13NO2	005255-69-6	15
4			6H-Benzo[d]pyrimido[5,4-b]thiopy...	215	C11H9N3S	1010267-44-6	15
5			3-(1,5-Dimethyl-1H-pyrazol-4-yl)...	215	C12H13N3O	1000275-23-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
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Sample : 03143-21
Misc :
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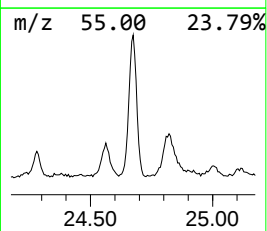
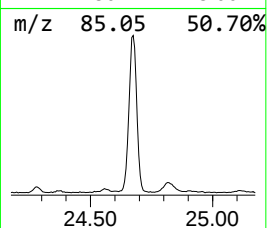
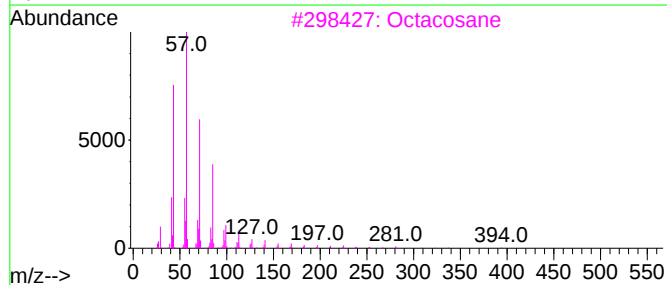
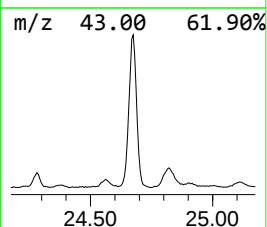
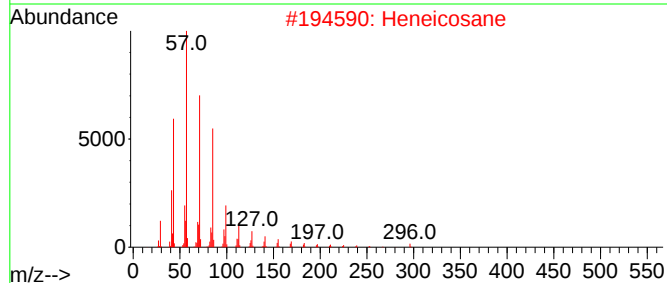
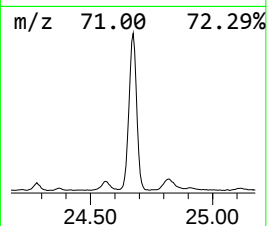
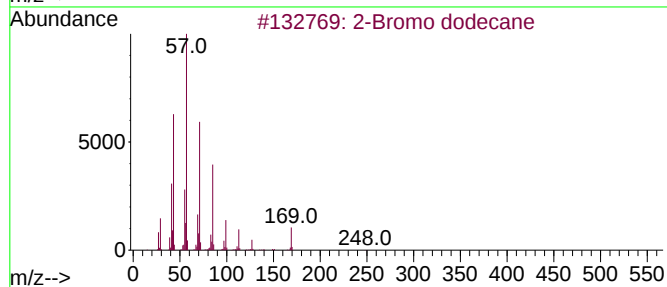
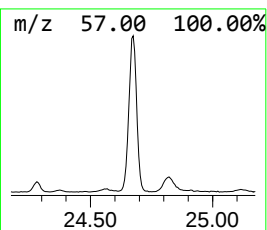
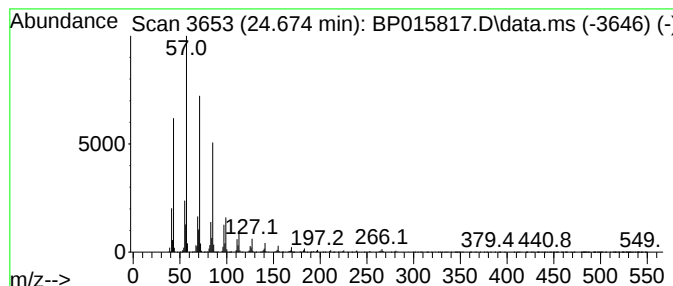
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-Bromo dodecane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.674	24.49 ng/ul	3699540	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromo dodecane			248	C12H25Br	013187-99-0	93
2	Heneicosane			296	C21H44	000629-94-7	91
3	Octacosane			394	C28H58	000630-02-4	91
4	Pentacosane			352	C25H52	000629-99-2	91
5	Hexatriacontane			507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
Operator : MA/JU
Sample : 03143-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFV8

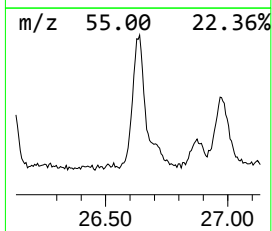
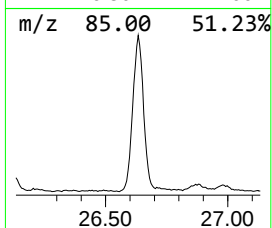
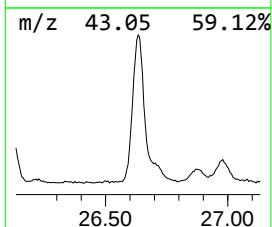
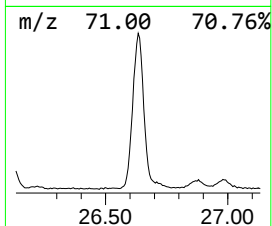
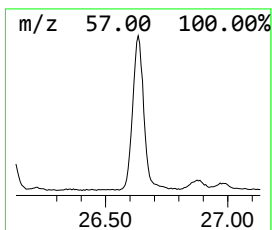
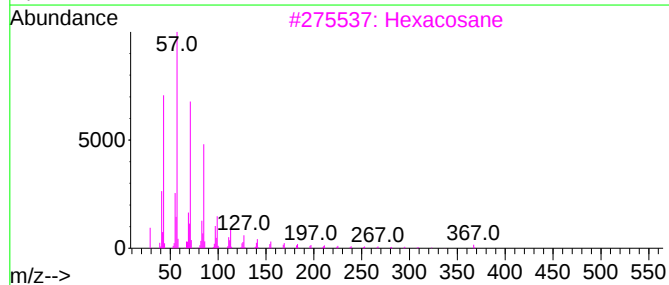
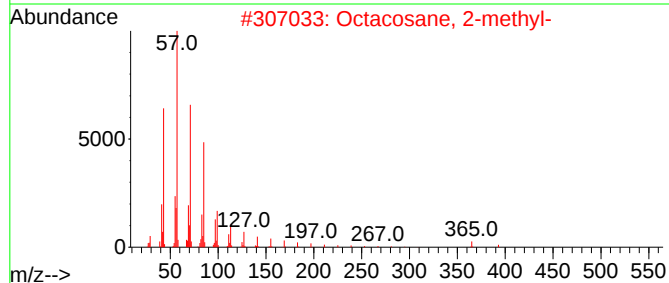
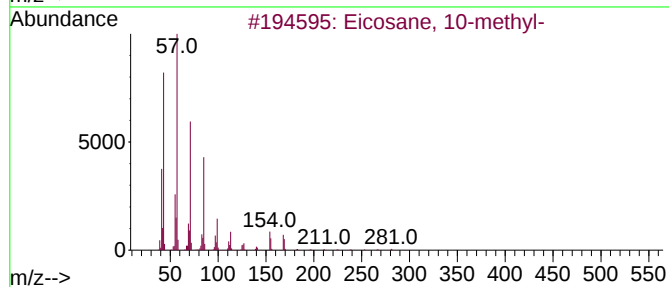
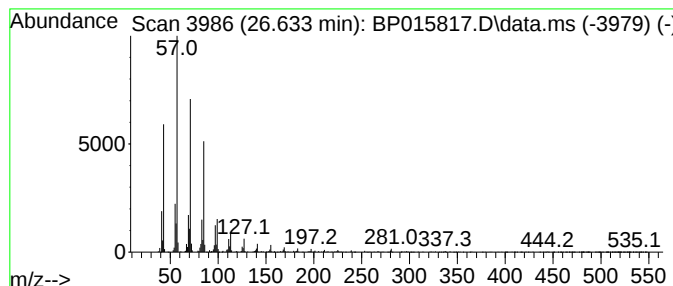
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Quant Title : SVOA CALIBRATION

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.
26.633	17.45 ng/ul	2636390	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	91
5		Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015817.D
Acq On : 29 Jun 2023 19:12
Operator : MA/JU
Sample : 03143-21
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFV8

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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	8.916	3.4	ng/ul	340894	1	8.058	1987240	20.0
n-Hexadecanoic ...	18.322	11.1	ng/ul	2143660	4	17.469	3857670	20.0
Octadecanoic acid	19.545	3.4	ng/ul	550237	5	21.557	3241770	20.0
(DEL) Alkane: S...	21.210	3.4	ng/ul	542600	5	21.557	3241770	20.0
Bis(2-ethylhexy...	21.416	2.2	ng/ul	360597	5	21.557	3241770	20.0
(DEL) Alkane: S...	22.133	2.3	ng/ul	372309	5	21.557	3241770	20.0
(DEL) Alkane: S...	23.239	10.0	ng/ul	1513020	6	24.098	3021180	20.0
unknown-01	24.563	7.5	ng/ul	1136650	6	24.098	3021180	20.0
2-Bromo dodecane	24.674	24.5	ng/ul	3699540	6	24.098	3021180	20.0
(DEL) Alkane: S...	26.633	17.4	ng/ul	2636390	6	24.098	3021180	20.0