

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
 Data File : BP015601.D
 Acq On : 17 Jun 2023 03:04
 Operator : MA/JU
 Sample : 03080-20
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFQ5

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: BP015601.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	33	37	52	rVB	41285	66616	1.28%	0.109%
2	3.711	85	89	96	rVB2	12250	17368	0.33%	0.028%
3	3.846	110	112	126	rBV	261887	456018	8.75%	0.743%
4	3.952	126	130	139	rVB2	41822	67571	1.30%	0.110%
5	4.075	147	151	156	rVB	19330	30535	0.59%	0.050%
6	4.175	165	168	174	rVB2	8850	12686	0.24%	0.021%
7	4.564	232	234	241	rVB8	6735	11535	0.22%	0.019%
8	4.758	264	267	272	rVB6	7198	9176	0.18%	0.015%
9	5.128	326	330	336	rVB4	4912	8039	0.15%	0.013%
10	6.299	523	529	535	rBV2	12466	25592	0.49%	0.042%
11	6.734	598	603	612	rVB	29458	50101	0.96%	0.082%
12	7.093	660	664	670	rVB9	5788	11803	0.23%	0.019%
13	7.240	682	689	702	rBV	750310	1264614	24.27%	2.061%
14	7.405	710	717	725	rBV	608142	942358	18.09%	1.536%
15	7.505	729	734	739	rVV5	10889	18128	0.35%	0.030%
16	7.610	745	752	761	rBV	797992	1304522	25.04%	2.126%
17	7.681	761	764	769	rVB6	9352	12725	0.24%	0.021%
18	8.081	823	832	844	rBV	576465	934651	17.94%	1.523%
19	8.799	947	954	969	rBV	839965	1434775	27.54%	2.339%
20	8.928	972	976	981	rVB7	6449	12158	0.23%	0.020%
21	9.257	1025	1032	1044	rBV	776682	1350465	25.92%	2.201%
22	9.410	1054	1058	1068	rVB6	15357	28242	0.54%	0.046%
23	9.640	1090	1097	1107	rBV6	7070	17131	0.33%	0.028%
24	9.987	1148	1156	1167	rBV	722232	1220218	23.42%	1.989%
25	10.216	1189	1195	1196	rBV6	7188	11615	0.22%	0.019%
26	10.228	1196	1197	1206	rVB9	7262	13881	0.27%	0.023%
27	10.528	1241	1248	1271	rBV	934970	1777862	34.13%	2.898%
28	10.910	1301	1313	1320	rBV	824841	1392975	26.74%	2.271%
29	10.957	1320	1321	1326	rVB2	9344	6627	0.13%	0.011%
30	11.051	1330	1337	1358	rVV	545547	1113642	21.38%	1.815%
31	11.710	1445	1449	1454	rBV5	10023	18915	0.36%	0.031%
32	11.916	1477	1484	1486	rBV8	5857	10657	0.20%	0.017%
33	12.404	1559	1567	1568	rBV5	5497	8516	0.16%	0.014%
34	12.492	1577	1582	1590	rVB	19002	32331	0.62%	0.053%
35	12.569	1590	1595	1598	rBV7	4979	7489	0.14%	0.012%
36	13.010	1666	1670	1675	rBV8	5497	9725	0.19%	0.016%

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Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	13.087	1678	1683	1689	rVV9	6958	14841	0.28%	0.024%
38	13.328	1719	1724	1731	rVB7	10407	15634	0.30%	0.025%
39	13.528	1755	1758	1762	rBV6	5043	7932	0.15%	0.013%
40	13.604	1766	1771	1777	rVB3	30218	55347	1.06%	0.090%
41	13.916	1821	1824	1829	rVB7	4319	7425	0.14%	0.012%
42	14.039	1840	1845	1852	rVB3	68300	98846	1.90%	0.161%
43	14.134	1852	1861	1868	rBV	1926943	2682183	51.49%	4.372%
44	14.239	1876	1879	1890	rVB	8451	17449	0.33%	0.028%
45	14.363	1896	1900	1903	rBV6	4177	7628	0.15%	0.012%
46	14.422	1903	1910	1926	rVV	2146453	3162205	60.70%	5.154%
47	14.598	1936	1940	1946	rVB6	9685	16505	0.32%	0.027%
48	14.675	1949	1953	1956	rBV4	22025	30152	0.58%	0.049%
49	14.728	1956	1962	1968	rVV	1286274	1867355	35.84%	3.044%
50	14.792	1969	1973	1981	rVB10	11900	26451	0.51%	0.043%
51	14.939	1990	1998	2019	rBV	501146	1240110	23.80%	2.021%
52	15.098	2021	2025	2027	rVV4	19582	33322	0.64%	0.054%
53	15.122	2027	2029	2037	rVB5	31707	54645	1.05%	0.089%
54	15.398	2073	2076	2081	rVB7	8390	11703	0.22%	0.019%
55	15.504	2088	2094	2097	rBV5	10617	19417	0.37%	0.032%
56	15.539	2097	2100	2104	rVB2	10462	17452	0.33%	0.028%
57	15.628	2112	2115	2119	rVB5	10553	12886	0.25%	0.021%
58	15.722	2124	2131	2137	rBV	2600950	3866675	74.22%	6.303%
59	15.845	2147	2152	2167	rBV	841382	1233351	23.67%	2.010%
60	15.957	2168	2171	2177	rVB4	21870	33993	0.65%	0.055%
61	16.081	2187	2192	2200	rBV	460970	634659	12.18%	1.034%
62	16.286	2223	2227	2233	rBV9	9685	16728	0.32%	0.027%
63	16.422	2246	2250	2251	rBV4	13817	17568	0.34%	0.029%
64	16.439	2251	2253	2258	rVB5	18366	21286	0.41%	0.035%
65	16.651	2285	2289	2294	rVB6	9651	14546	0.28%	0.024%
66	16.863	2321	2325	2330	rBV5	19036	24565	0.47%	0.040%
67	17.010	2347	2350	2356	rVB3	17923	25090	0.48%	0.041%
68	17.216	2381	2385	2389	rBV5	16150	22654	0.43%	0.037%
69	17.357	2405	2409	2416	rBV4	37736	63626	1.22%	0.104%
70	17.422	2417	2420	2424	rBV6	15352	20326	0.39%	0.033%
71	17.480	2424	2430	2436	rBV	1572842	2300203	44.15%	3.749%
72	17.528	2436	2438	2441	rVV	35527	39412	0.76%	0.064%
73	17.580	2441	2447	2461	rVB2	2764134	4119507	79.07%	6.715%
74	17.757	2473	2477	2480	rBV	9500	17070	0.33%	0.028%
75	17.822	2486	2488	2490	rBV2	8398	6456	0.12%	0.011%
76	18.127	2537	2540	2544	rVB5	11804	14361	0.28%	0.023%

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77	18.227	2554	2557	2562	rVB6	24271	33866	0.65%	0.055%
78	18.286	2563	2567	2570	rBV2	35498	45806	0.88%	0.075%
79	18.345	2572	2577	2586	rBV	1277203	1652342	31.72%	2.693%
80	18.622	2620	2624	2630	rBV5	28676	65156	1.25%	0.106%
81	19.222	2724	2726	2732	rVB3	37926	45261	0.87%	0.074%
82	19.380	2750	2753	2757	rVB	49003	58714	1.13%	0.096%
83	19.427	2758	2761	2763	rBV3	50000	63062	1.21%	0.103%
84	19.451	2763	2765	2767	rVV2	98146	111075	2.13%	0.181%
85	19.480	2767	2770	2777	rVB	134297	182847	3.51%	0.298%
86	19.569	2780	2785	2792	rBV	439523	628300	12.06%	1.024%
87	19.810	2820	2826	2835	rBV	3785831	5175774	99.35%	8.437%
88	20.780	2988	2991	2993	rBV2	91212	101766	1.95%	0.166%
89	21.098	3041	3045	3050	rVB	275365	334080	6.41%	0.545%
90	21.245	3065	3070	3083	rVB2	794274	1312730	25.20%	2.140%
91	21.445	3101	3104	3109	rVB	115684	131331	2.52%	0.214%
92	21.563	3119	3124	3129	rBV2	1784909	2495538	47.90%	4.068%
93	22.163	3224	3226	3230	rVV	203981	219394	4.21%	0.358%
94	22.204	3230	3233	3242	rVB5	218435	391640	7.52%	0.638%
95	22.692	3313	3316	3321	rVB	230036	304350	5.84%	0.496%
96	23.280	3411	3416	3422	rBV	787795	1250855	24.01%	2.039%
97	23.951	3523	3530	3543	rVB2	2508753	5209636	100.00%	8.492%
98	24.110	3551	3557	3566	rVB	1078492	2268099	43.54%	3.697%
99	24.733	3657	3663	3672	rVB	927283	2046002	39.27%	3.335%
100	24.857	3678	3684	3699	rVB4	494678	1619016	31.08%	2.639%

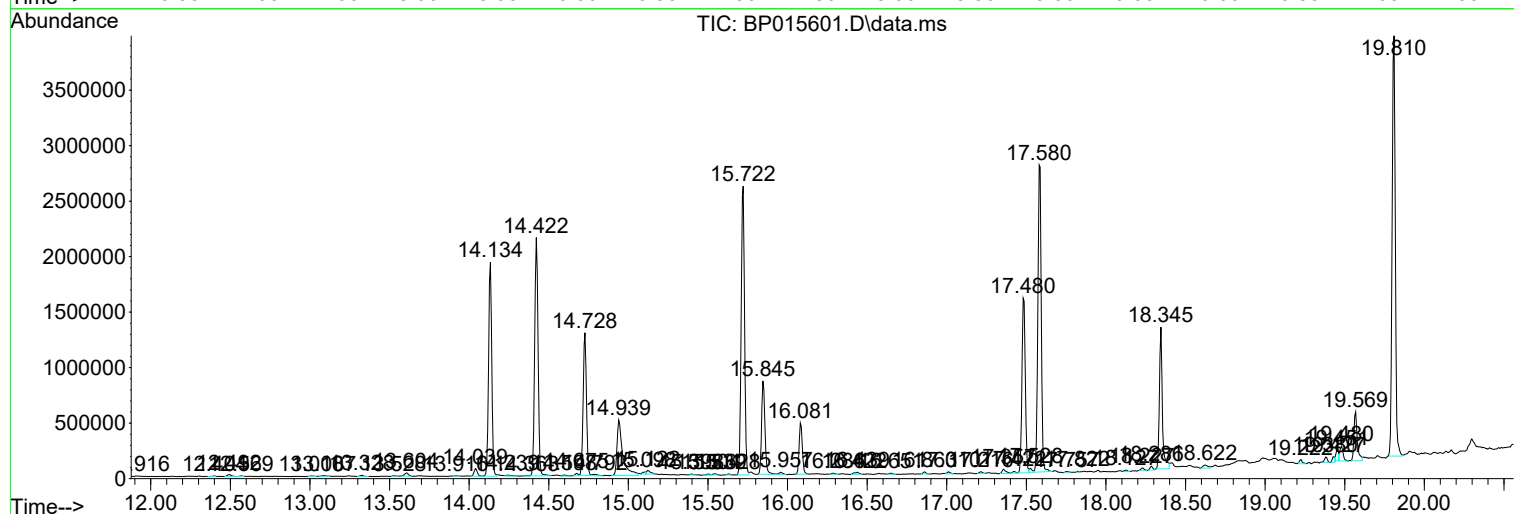
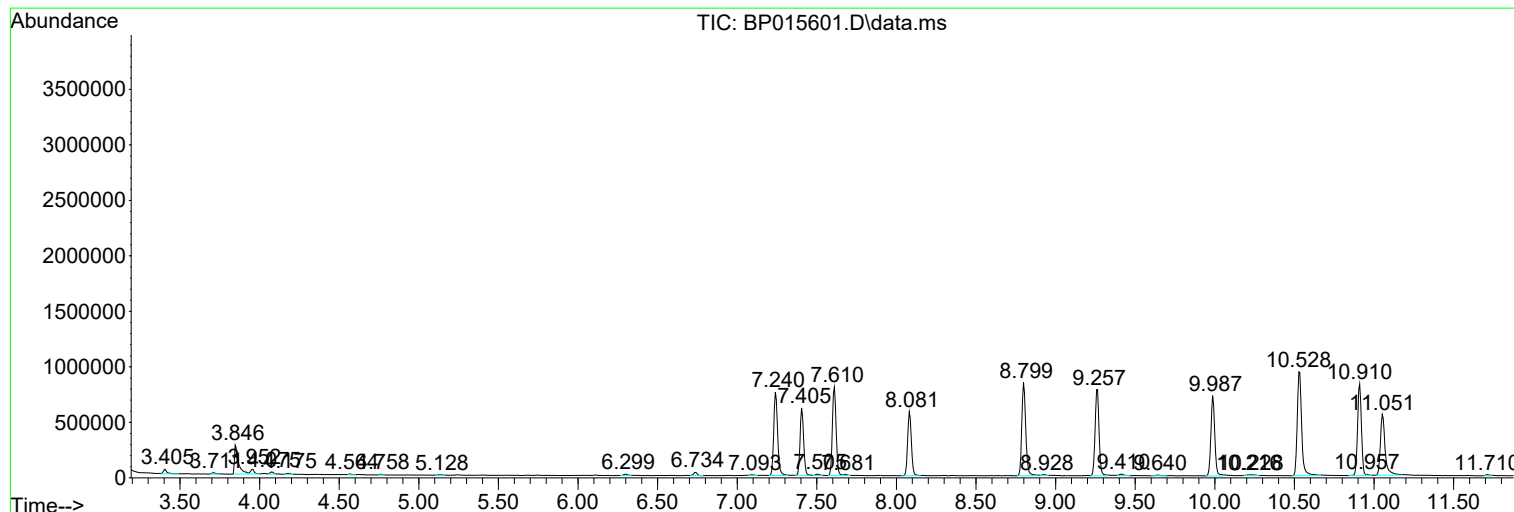
Sum of corrected areas: 61349466

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

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



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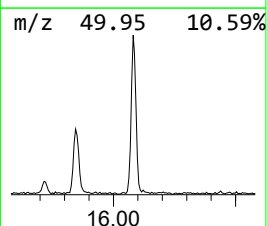
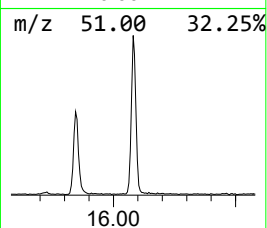
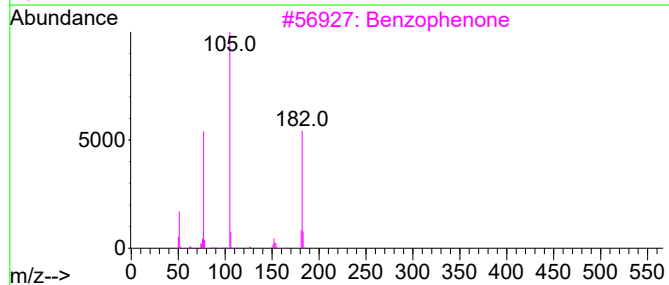
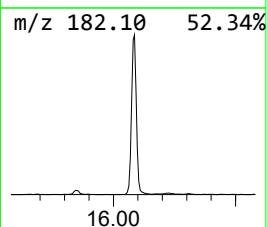
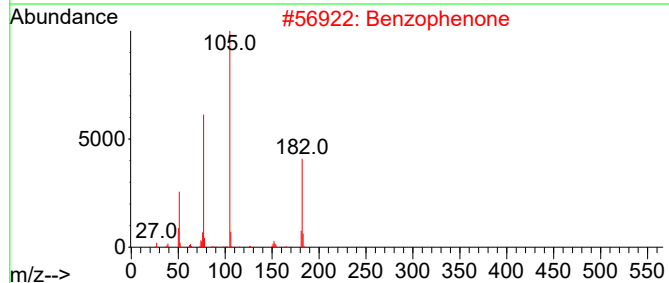
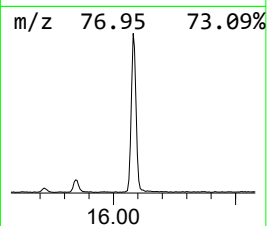
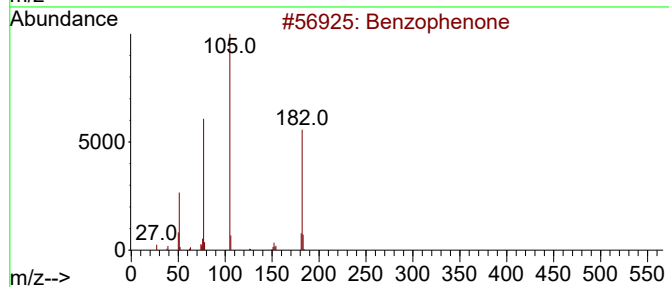
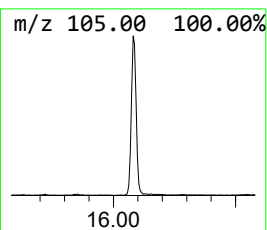
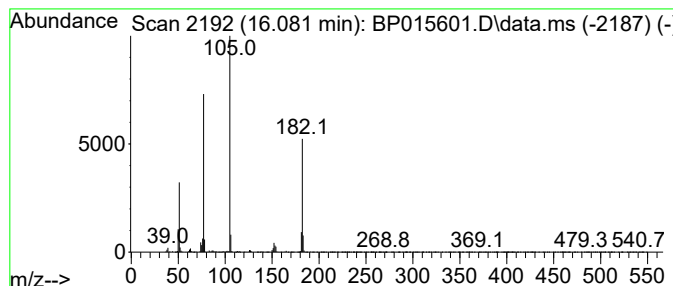
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.081	6.80 ng/ul	634659	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	95
5			Benzophenone	182	C13H10O	000119-61-9	91



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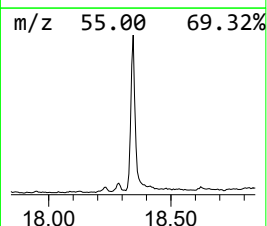
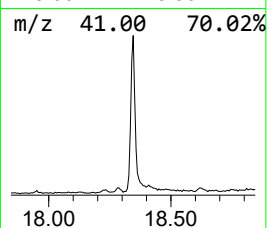
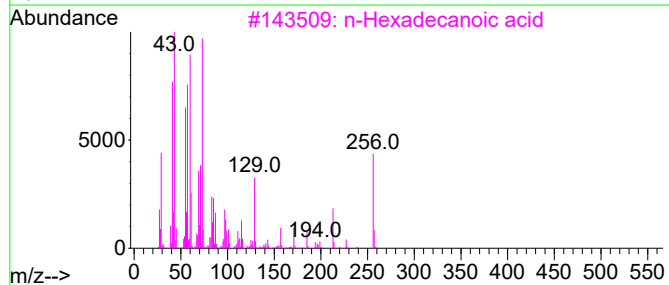
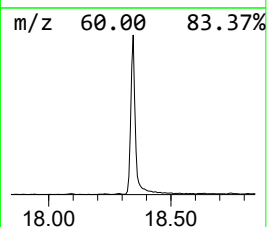
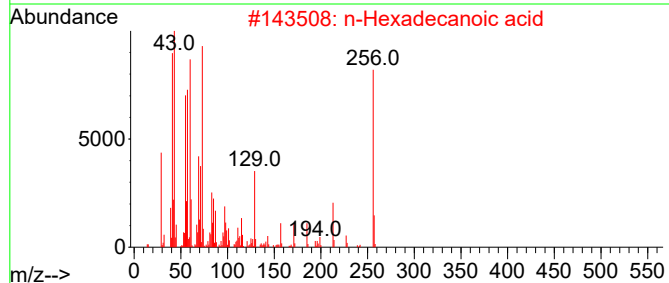
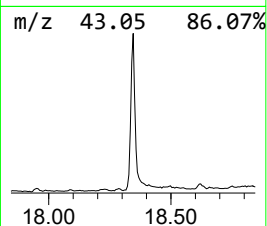
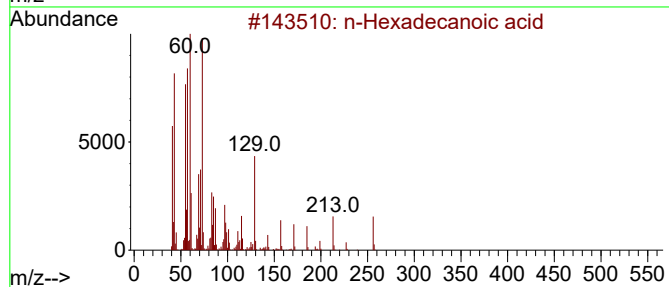
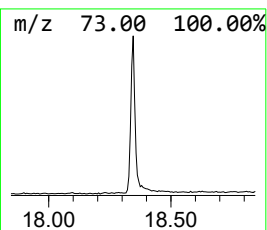
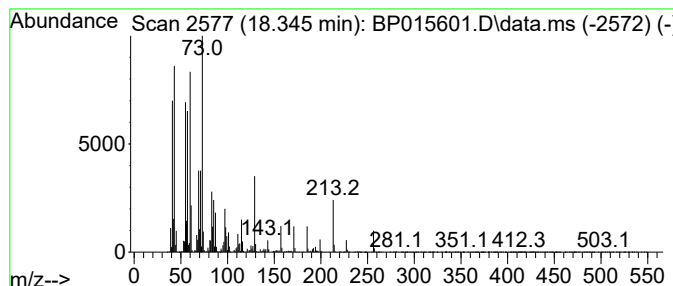
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	14.37 ng/ul	1652340	Phenanthrene-d10	17.480

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97



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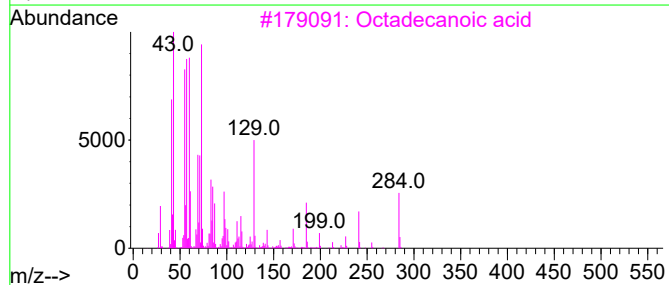
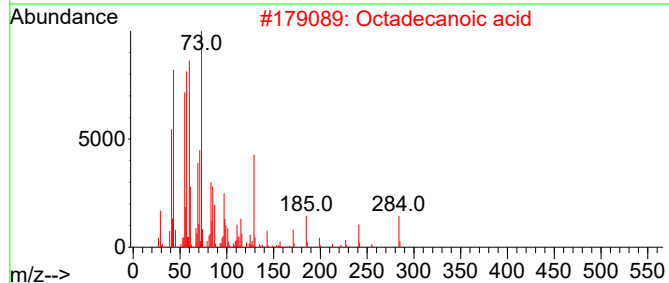
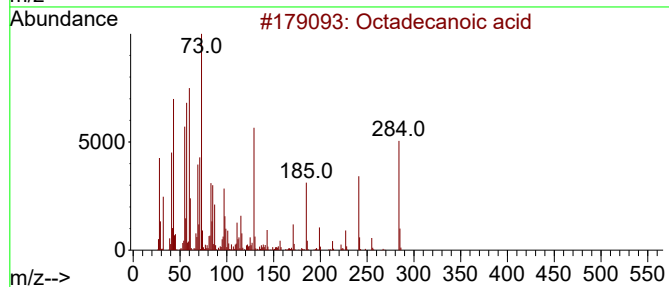
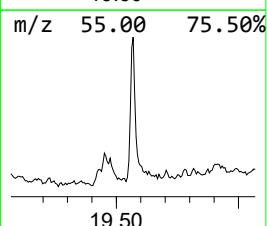
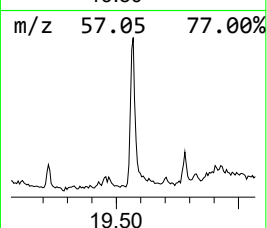
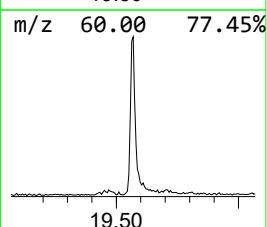
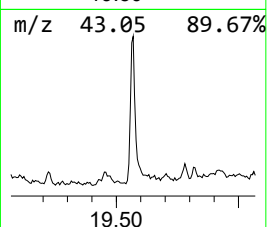
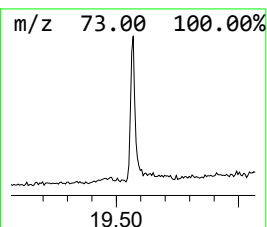
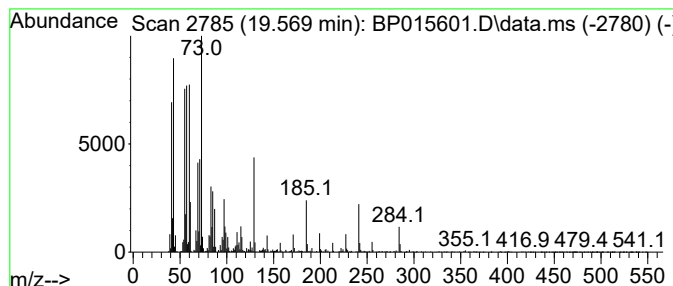
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Quant Title : SVOA CALIBRATION

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.
19.569	5.04 ng/ul	628300	Chrysene-d12	21.563

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			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	91



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Operator : MA/JU
Sample : 03080-20
Misc :
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ClientSampleId :
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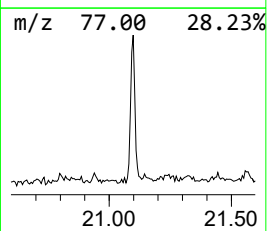
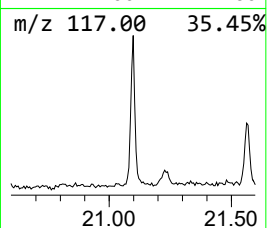
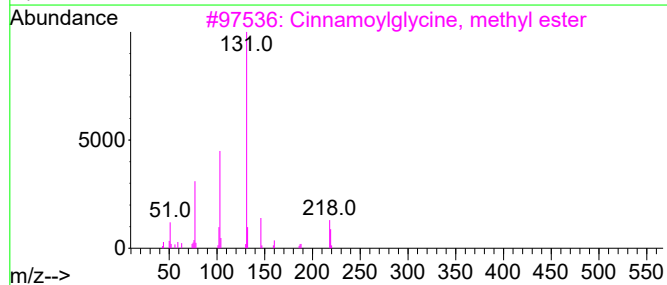
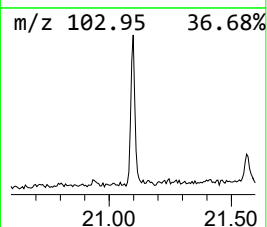
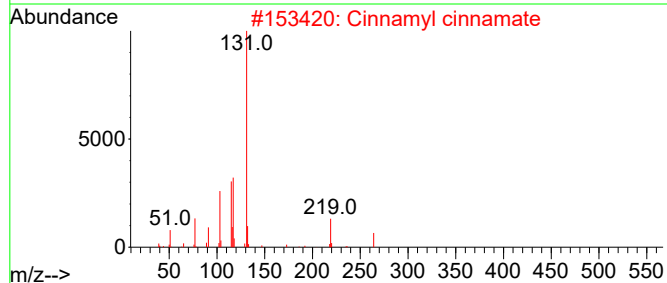
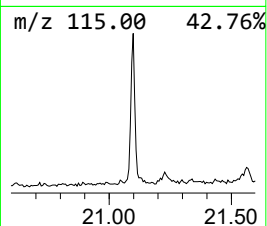
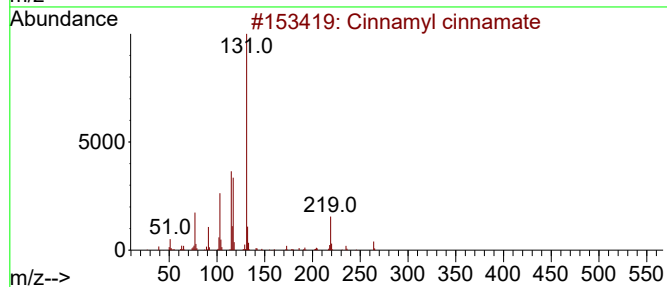
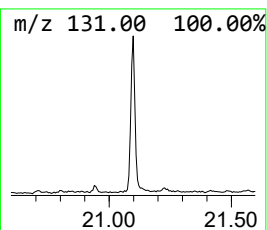
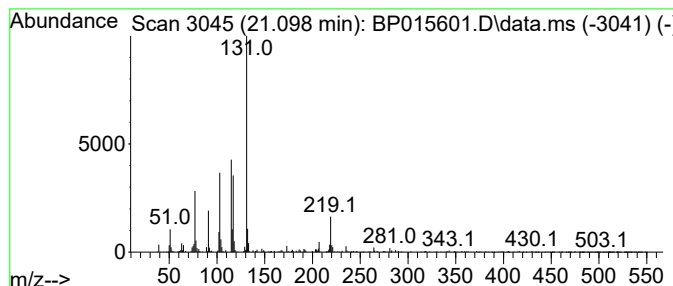
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cinnamyl cinnamate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.098	2.68 ng/ul	334080	Chrysene-d12	21.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	91
2			Cinnamyl cinnamate	264	C18H16O2	000122-69-0	86
3			Cinnamoylglycine, methyl ester	219	C12H13NO3	040778-04-9	50
4			(E)-4-Bromo-4-methyl-1-phenylpen...	252	C12H13BrO	1000443-02-8	50
5			N-(3-Chlorophenyl)-3-phenyl-2-pr...	257	C15H12ClNO	064741-15-7	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
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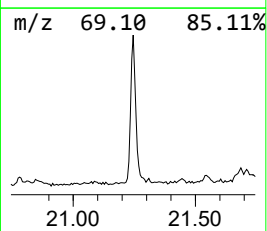
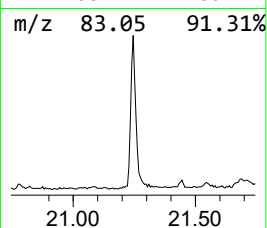
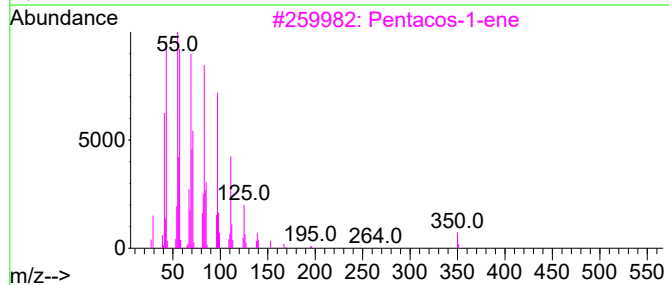
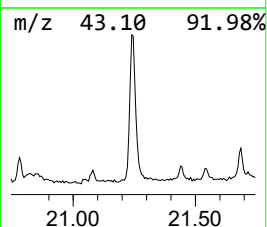
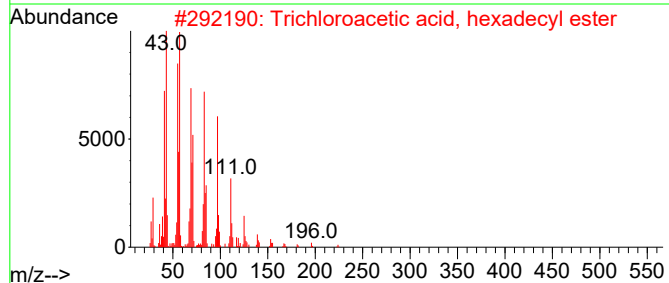
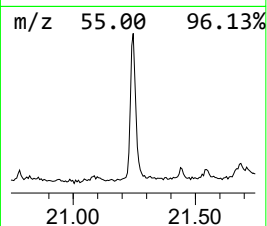
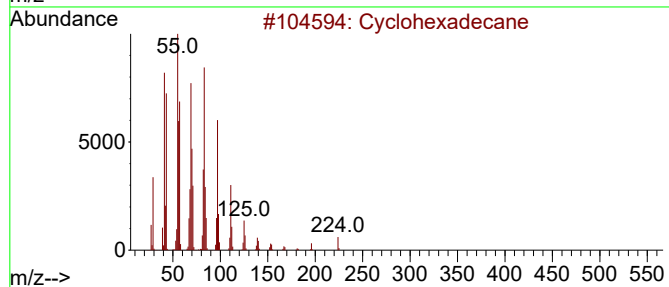
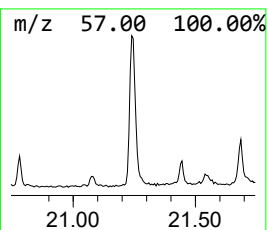
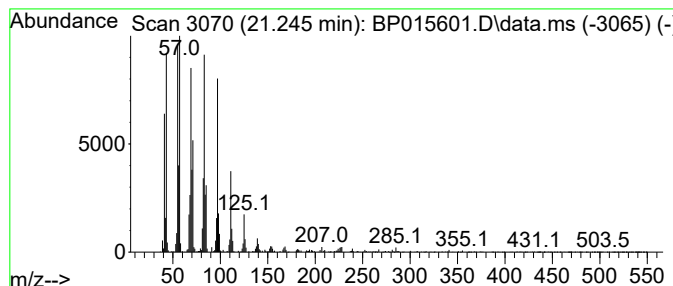
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic21.245 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.245	10.52 ng/ul	1312730	Chrysene-d12	21.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
5			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015601.D
Acq On : 17 Jun 2023 03:04
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Sample : 03080-20
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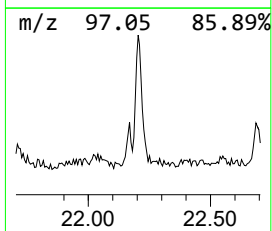
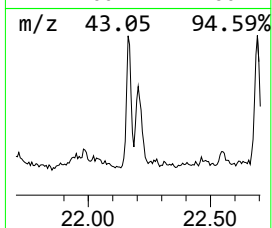
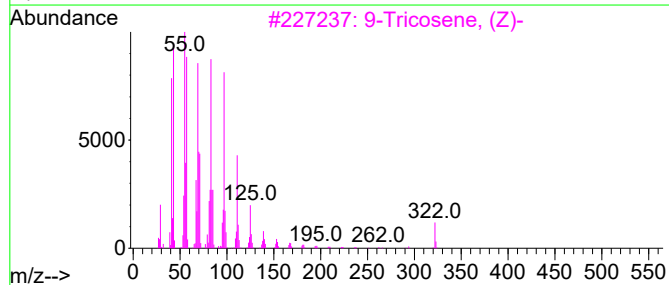
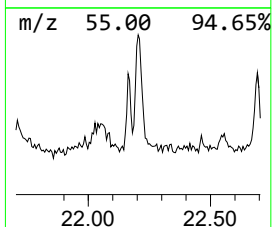
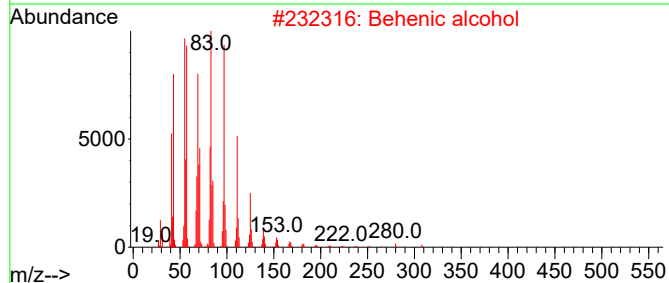
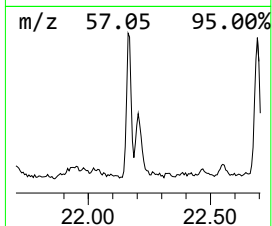
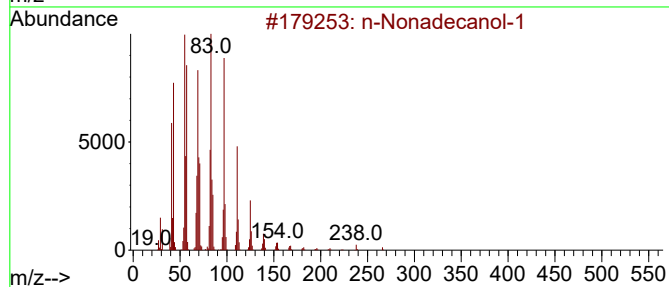
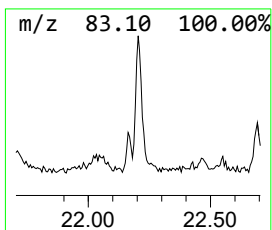
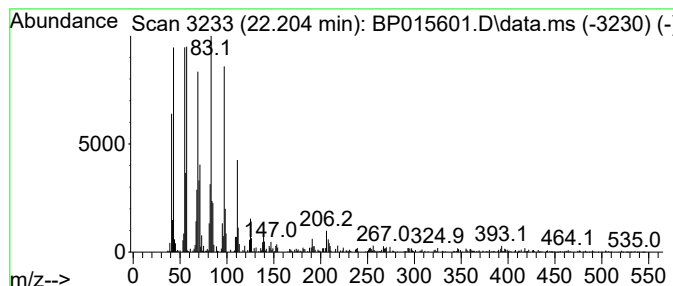
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Nonadecanol-1 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.204	3.14 ng/ul	391640	Chrysene-d12	21.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Nonadecanol-1	284	C19H40O	001454-84-8	91
2			Behenic alcohol	326	C22H46O	000661-19-8	91
3			9-Tricosene, (Z)-	322	C23H46	027519-02-4	91
4			Bromoacetic acid, octadecyl ester	390	C20H39BrO2	018992-03-5	87
5			Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015601.D
Acq On : 17 Jun 2023 03:04
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Sample : 03080-20
Misc :
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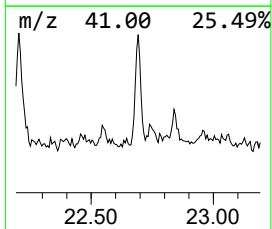
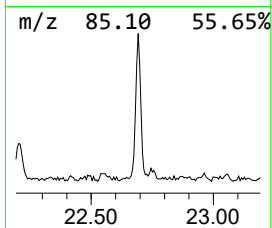
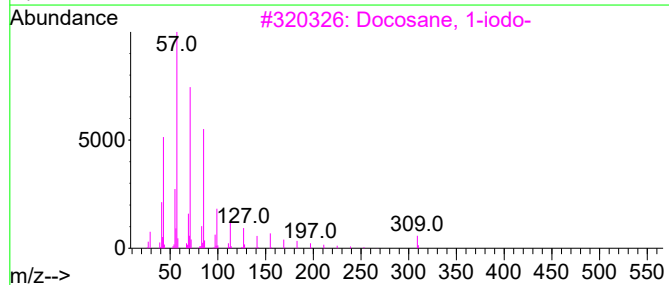
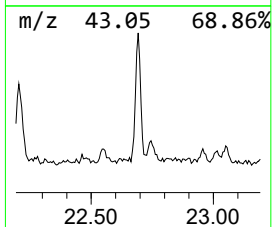
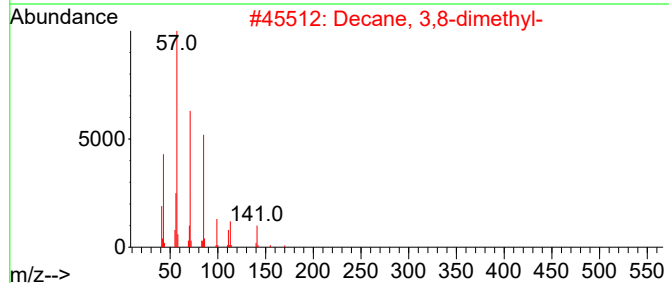
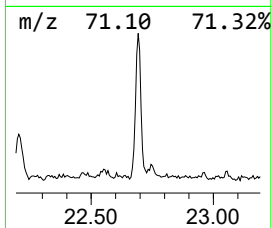
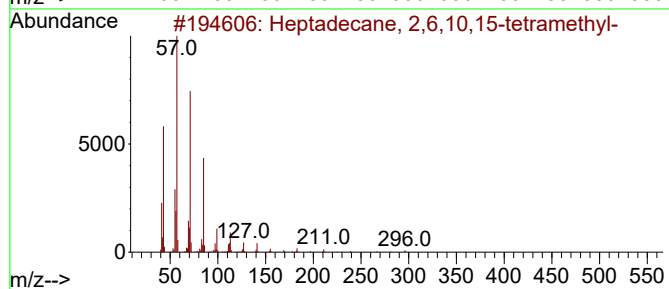
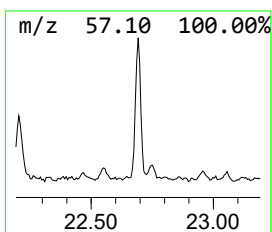
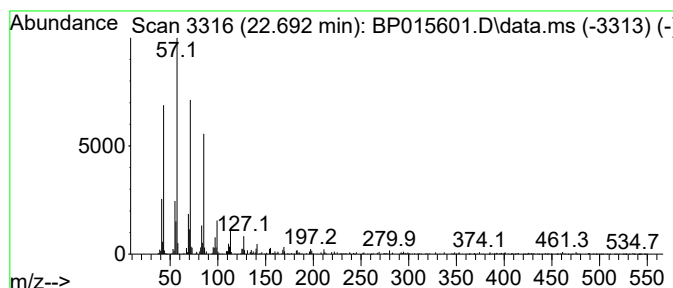
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.692	2.44 ng/ul	304350	Chrysene-d12	21.563	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	94
2	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	90
3	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	90
4	Octadecane, 5,14-dibutyl-	366	C26H54	055282-13-8	87
5	Tetracosane	338	C24H50	000646-31-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015601.D
Acq On : 17 Jun 2023 03:04
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Sample : 03080-20
Misc :
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Instrument :
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ClientSampleId :
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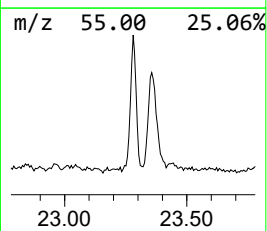
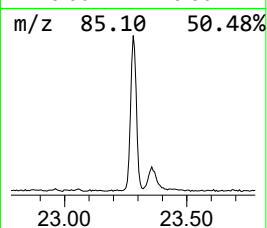
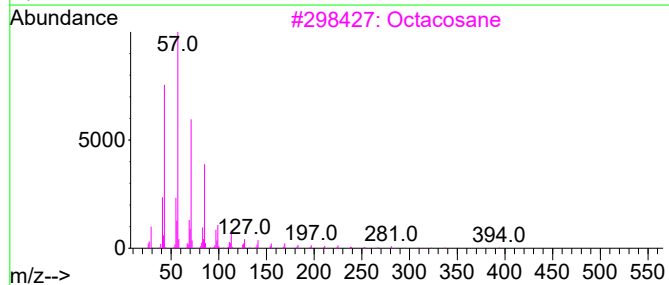
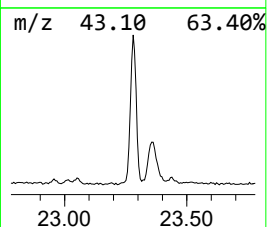
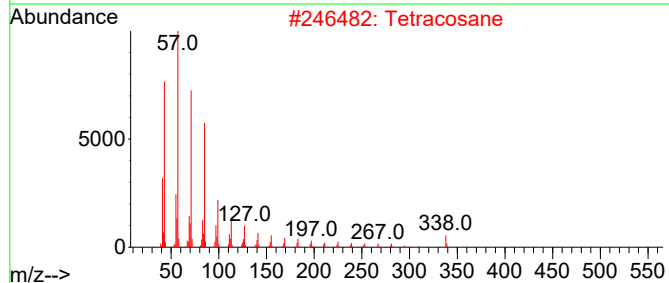
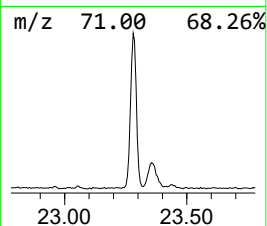
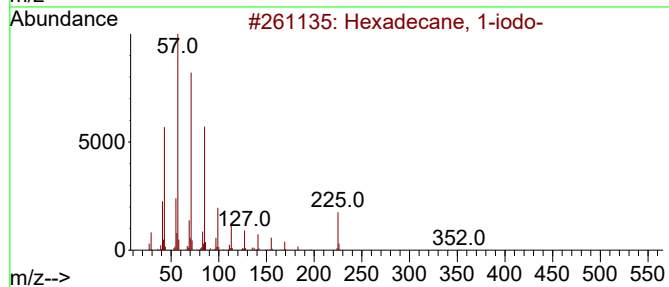
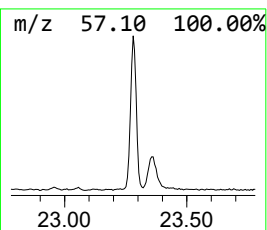
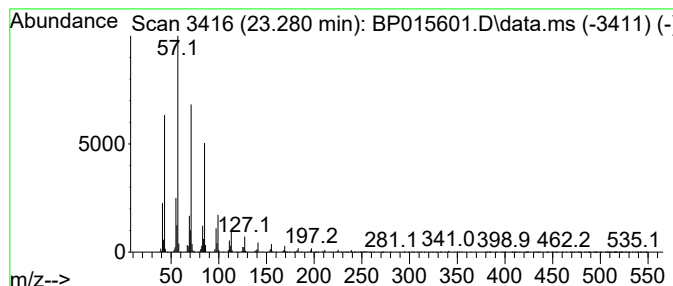
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Hexadecane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.280	11.03 ng/ul	1250860	Perylene-d12	24.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	95
2			Tetracosane	338	C24H50	000646-31-1	94
3			Octacosane	394	C28H58	000630-02-4	91
4			Tetracosane	338	C24H50	000646-31-1	91
5			Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
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Acq On : 17 Jun 2023 03:04
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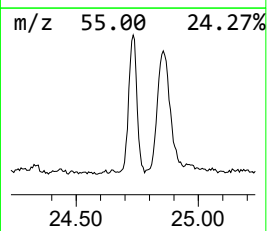
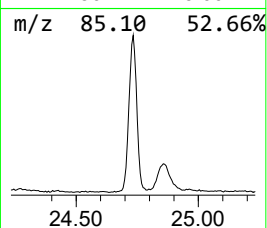
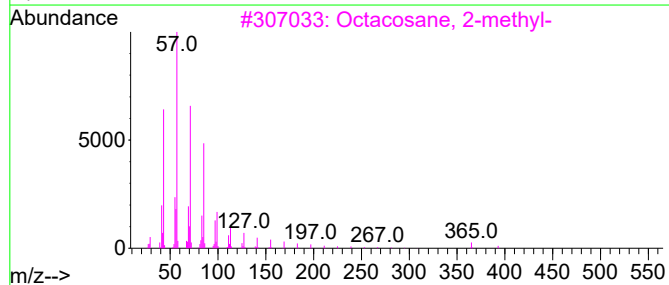
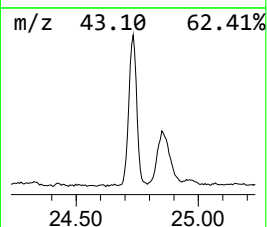
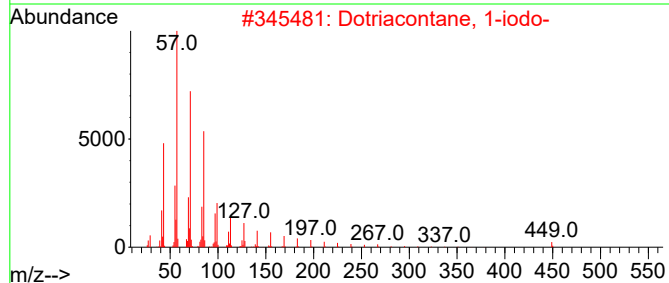
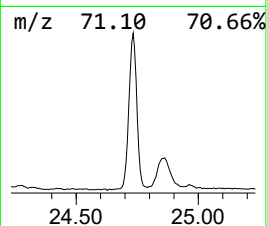
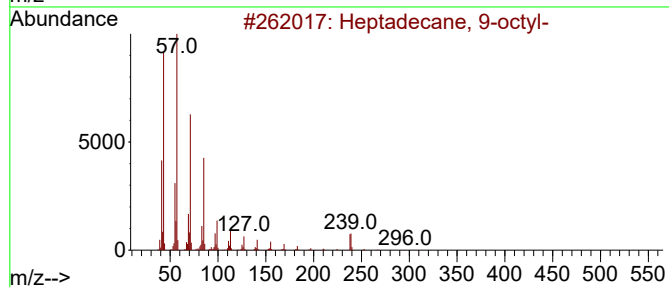
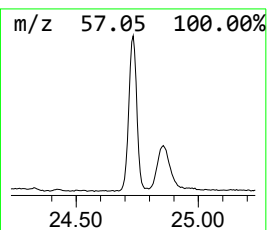
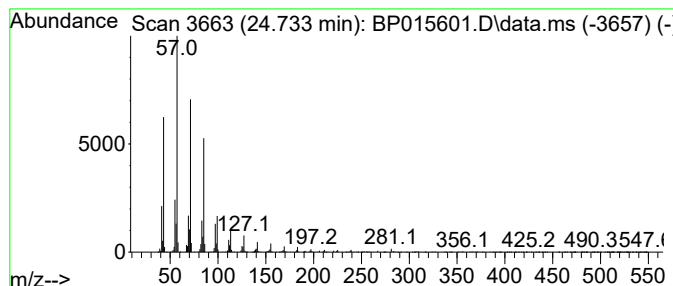
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.733	18.04 ng/ul	2046000	Perylene-d12	24.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Octacosane	394	C28H58	000630-02-4	91
5			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015601.D
Acq On : 17 Jun 2023 03:04
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ALS Vial : 32 Sample Multiplier: 1

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ClientSampleId :
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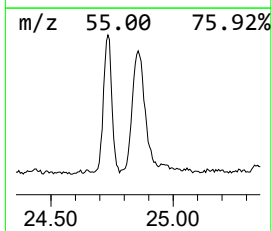
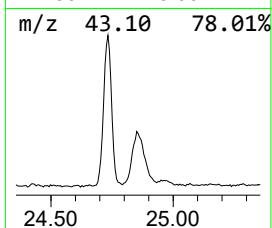
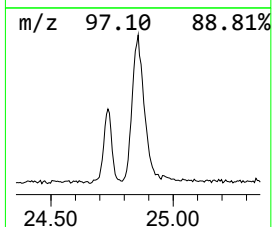
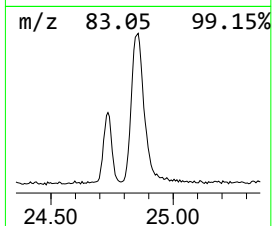
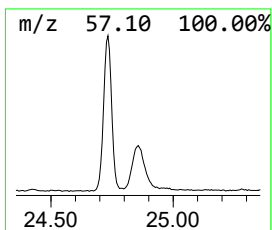
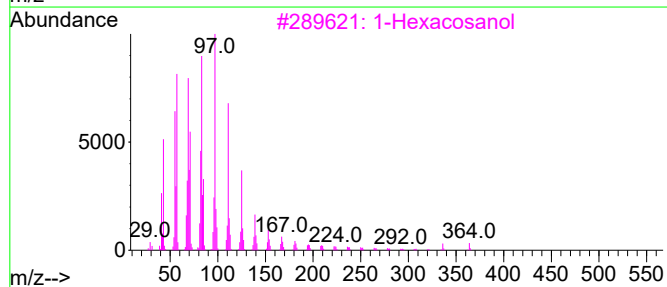
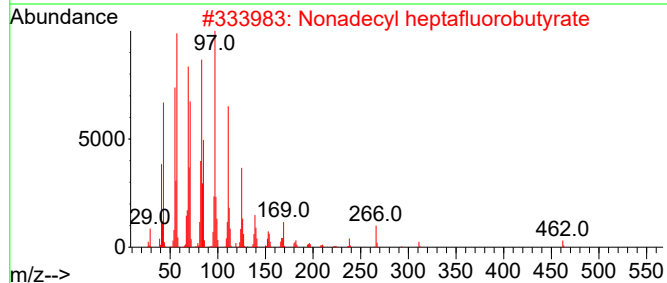
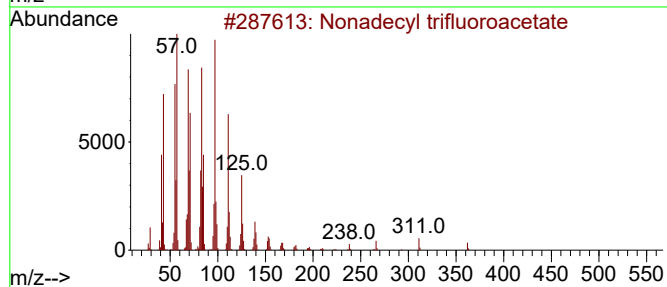
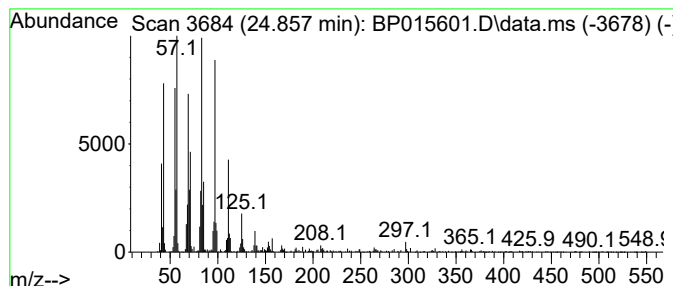
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Nonadecyl trifluoroacetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.857	14.28 ng/ul	1619020	Perylene-d12	24.110

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	76
2			Nonadecyl heptafluorobutyrate	480	C23H39F7O2	1000351-83-9	76
3			1-Hexacosanol	382	C26H54O	000506-52-5	76
4			1-Hexacosene	364	C26H52	018835-33-1	70
5			1-Docosene	308	C22H44	001599-67-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061623\
Data File : BP015601.D
Acq On : 17 Jun 2023 03:04
Operator : MA/JU
Sample : 03080-20
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFQ5

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
Benzophenone	16.081	6.8	ng/ul	634659	3	14.728	1867360	20.0
n-Hexadecanoic ...	18.345	14.4	ng/ul	1652340	4	17.480	2300200	20.0
Octadecanoic acid	19.569	5.0	ng/ul	628300	5	21.563	2495540	20.0
Cinnamyl cinnamate	21.098	2.7	ng/ul	334080	5	21.563	2495540	20.0
(DEL) Alkane: C...	21.245	10.5	ng/ul	1312730	5	21.563	2495540	20.0
n-Nonadecanol-1	22.204	3.1	ng/ul	391640	5	21.563	2495540	20.0
(DEL) Alkane: S...	22.692	2.4	ng/ul	304350	5	21.563	2495540	20.0
Hexadecane, 1-i...	23.280	11.0	ng/ul	1250860	6	24.110	2268100	20.0
(DEL) Alkane: S...	24.733	18.0	ng/ul	2046000	6	24.110	2268100	20.0
Nonadecyl trifl...	24.857	14.3	ng/ul	1619020	6	24.110	2268100	20.0