

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015863.D
 Acq On : 01 Jul 2023 01:54
 Operator : MA/JU
 Sample : 03161-12
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFW8

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: BP015863.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	28	32	40	rVB	56450	83084	1.25%	0.069%
2	3.822	106	108	121	rVV	478340	707317	10.67%	0.590%
3	3.922	121	125	135	rVB	299836	400583	6.04%	0.334%
4	3.999	135	138	141	rBV	34082	41019	0.62%	0.034%
5	4.046	141	146	150	rVB	45189	66804	1.01%	0.056%
6	4.146	160	163	172	rVB2	22460	41897	0.63%	0.035%
7	4.722	257	261	270	rVB2	53316	86422	1.30%	0.072%
8	5.216	342	345	353	rBV2	25068	49001	0.74%	0.041%
9	7.234	681	688	702	rBV	562464	1470948	22.19%	1.227%
10	7.375	706	712	725	rVV	693421	1192950	18.00%	0.995%
11	7.581	741	747	766	rVV	838503	1579012	23.82%	1.317%
12	8.052	819	827	841	rVV	859794	1585949	23.93%	1.322%
13	8.346	872	877	888	rVB2	44151	79090	1.19%	0.066%
14	8.793	946	953	971	rBV	424842	1240958	18.72%	1.035%
15	9.240	1022	1029	1051	rBV	716892	1558007	23.51%	1.299%
16	9.969	1144	1153	1179	rBV	503976	1338749	20.20%	1.116%
17	10.546	1243	1251	1281	rBV	537713	1913146	28.87%	1.595%
18	10.881	1301	1308	1326	rVV	1172058	2385879	36.00%	1.989%
19	11.081	1335	1342	1373	rVB2	110289	466518	7.04%	0.389%
20	12.757	1621	1627	1633	rBV7	17040	40337	0.61%	0.034%
21	13.022	1665	1672	1678	rBV9	13252	39718	0.60%	0.033%
22	13.498	1748	1753	1758	rVB2	23411	35742	0.54%	0.030%
23	13.581	1758	1767	1774	rBV2	31071	66026	1.00%	0.055%
24	14.016	1835	1841	1847	rBV5	17025	42237	0.64%	0.035%
25	14.110	1847	1857	1870	rBV	1774905	2931845	44.24%	2.445%
26	14.398	1896	1906	1922	rBV	2162423	3519937	53.11%	2.935%
27	14.569	1930	1935	1940	rVB2	30743	46126	0.70%	0.038%
28	14.704	1951	1958	1977	rVB	1852088	3005894	45.35%	2.506%
29	14.910	1985	1993	2000	rBV2	106475	243537	3.67%	0.203%
30	14.987	2000	2006	2022	rVV2	213165	744948	11.24%	0.621%
31	15.116	2024	2028	2036	rVB8	40752	101119	1.53%	0.084%
32	15.469	2083	2088	2093	rBV6	18009	35008	0.53%	0.029%
33	15.516	2093	2096	2101	rVB	39899	53489	0.81%	0.045%
34	15.698	2120	2127	2146	rBV	2399247	4150831	62.63%	3.461%
35	15.863	2150	2155	2175	rBV	402110	1020163	15.39%	0.851%
36	16.069	2184	2190	2199	rBV	620826	993609	14.99%	0.828%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	16.398	2239	2246	2253	rBV3	55077	126760	1.91%	0.106%
38	16.545	2267	2271	2277	rBV4	60711	101025	1.52%	0.084%
39	16.610	2277	2282	2294	rVB2	154056	304692	4.60%	0.254%
40	16.839	2318	2321	2327	rBV4	37741	48247	0.73%	0.040%
41	17.181	2375	2379	2382	rBV3	49058	57196	0.86%	0.048%
42	17.334	2400	2405	2412	rBV	229595	359535	5.42%	0.300%
43	17.398	2412	2416	2421	rVV	151780	220097	3.32%	0.184%
44	17.463	2421	2427	2432	rVV	1897635	2930292	44.21%	2.443%
45	17.504	2432	2434	2439	rVV	225939	324289	4.89%	0.270%
46	17.563	2439	2444	2457	rVV2	1931841	3380400	51.00%	2.819%
47	17.916	2500	2504	2511	rVB3	65470	96707	1.46%	0.081%
48	18.063	2526	2529	2532	rBV3	42157	51292	0.77%	0.043%
49	18.092	2532	2534	2539	rVB5	44823	49751	0.75%	0.041%
50	18.145	2539	2543	2545	rBV5	30896	47623	0.72%	0.040%
51	18.204	2548	2553	2557	rBV2	159594	257183	3.88%	0.214%
52	18.263	2558	2563	2568	rBV	1030837	1316118	19.86%	1.097%
53	18.316	2568	2572	2580	rBV	3055960	4093531	61.76%	3.413%
54	18.604	2616	2621	2625	rBV5	146254	266504	4.02%	0.222%
55	19.039	2692	2695	2701	rBV8	62871	100131	1.51%	0.083%
56	19.104	2702	2706	2712	rBV8	93572	211593	3.19%	0.176%
57	19.192	2718	2721	2728	rVB4	242675	419137	6.32%	0.349%
58	19.298	2736	2739	2743	rVB2	76195	92331	1.39%	0.077%
59	19.369	2747	2751	2753	rBV2	142872	204988	3.09%	0.171%
60	19.404	2753	2757	2759	rVV	346340	440229	6.64%	0.367%
61	19.428	2759	2761	2762	rVV	234599	223455	3.37%	0.186%
62	19.463	2762	2767	2776	rVB2	697814	1387515	20.93%	1.157%
63	19.539	2776	2780	2784	rBV	954016	1229029	18.54%	1.025%
64	19.757	2812	2817	2818	rBV3	333346	475604	7.18%	0.397%
65	19.792	2818	2823	2832	rVB2	2827503	4690226	70.77%	3.911%
66	19.933	2844	2847	2851	rVB3	496293	618904	9.34%	0.516%
67	19.986	2853	2856	2859	rVB4	195683	228490	3.45%	0.191%
68	20.027	2860	2863	2864	rBV2	128009	144635	2.18%	0.121%
69	20.098	2871	2875	2878	rBV5	181426	339501	5.12%	0.283%
70	20.133	2878	2881	2885	rVV3	512291	625689	9.44%	0.522%
71	20.175	2885	2888	2892	rVV2	225825	274259	4.14%	0.229%
72	20.227	2893	2897	2900	rVB4	417579	535570	8.08%	0.447%
73	20.269	2900	2904	2912	rBV3	481374	996115	15.03%	0.831%
74	20.339	2912	2916	2923	rVB4	601199	976217	14.73%	0.814%
75	20.410	2924	2928	2931	rBV2	465131	609522	9.20%	0.508%
76	20.457	2932	2936	2940	rBV2	716446	984361	14.85%	0.821%

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77	20.575	2951	2956	2962	rBV6	1002408	1691404	25.52%	1.410%
78	20.633	2962	2966	2969	rVV3	435065	566320	8.54%	0.472%
79	20.827	2996	2999	3003	rBV2	454323	532995	8.04%	0.444%
80	20.874	3004	3007	3020	rVB5	1058632	2066669	31.18%	1.723%
81	20.998	3022	3028	3034	rBV6	726928	1537423	23.20%	1.282%
82	21.051	3034	3037	3043	rVV3	370495	503431	7.60%	0.420%
83	21.163	3053	3056	3059	rBV2	575206	650145	9.81%	0.542%
84	21.210	3059	3064	3071	rVV4	1843860	3246295	48.98%	2.707%
85	21.274	3072	3075	3081	rVB5	618525	923522	13.93%	0.770%
86	21.398	3091	3096	3102	rBV2	1369847	2449027	36.95%	2.042%
87	21.551	3118	3122	3128	rBV2	1603374	2251449	33.97%	1.877%
88	21.898	3178	3181	3185	rBV4	394302	478360	7.22%	0.399%
89	22.127	3217	3220	3223	rBV	528569	601140	9.07%	0.501%
90	22.169	3223	3227	3241	rVB	1348737	2410547	36.37%	2.010%
91	22.916	3351	3354	3360	rVB	378873	508419	7.67%	0.424%
92	23.233	3402	3408	3414	rVB	3106662	5098567	76.93%	4.251%
93	23.316	3417	3422	3450	rVB	836295	2684541	40.50%	2.238%
94	23.927	3521	3526	3532	rVB2	1436097	2925792	44.14%	2.440%
95	24.092	3549	3554	3571	rVB	1290858	3211498	48.46%	2.678%
96	24.274	3581	3585	3592	rVB	669495	1205671	18.19%	1.005%
97	24.557	3626	3633	3642	rVB2	3033739	6627759	100.00%	5.526%
98	24.663	3645	3651	3664	rVB	1667168	3798287	57.31%	3.167%
99	26.115	3890	3898	3911	rVB	1829149	4981095	75.16%	4.153%
100	26.621	3977	3984	4006	rVB	1705960	5783091	87.26%	4.822%

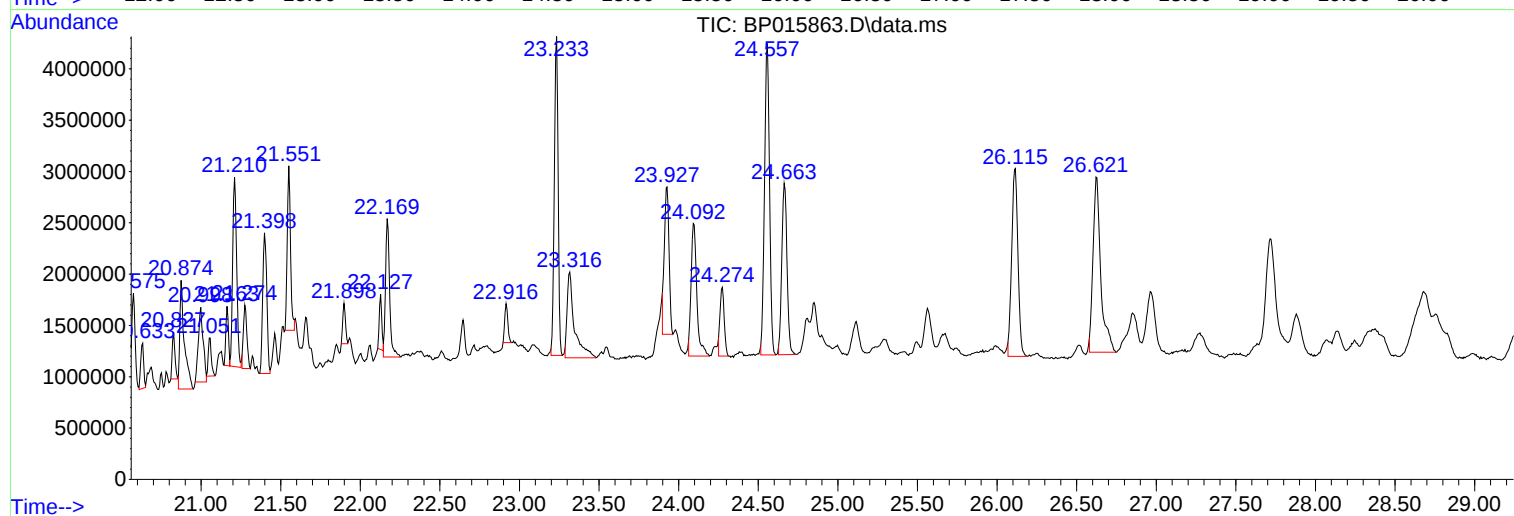
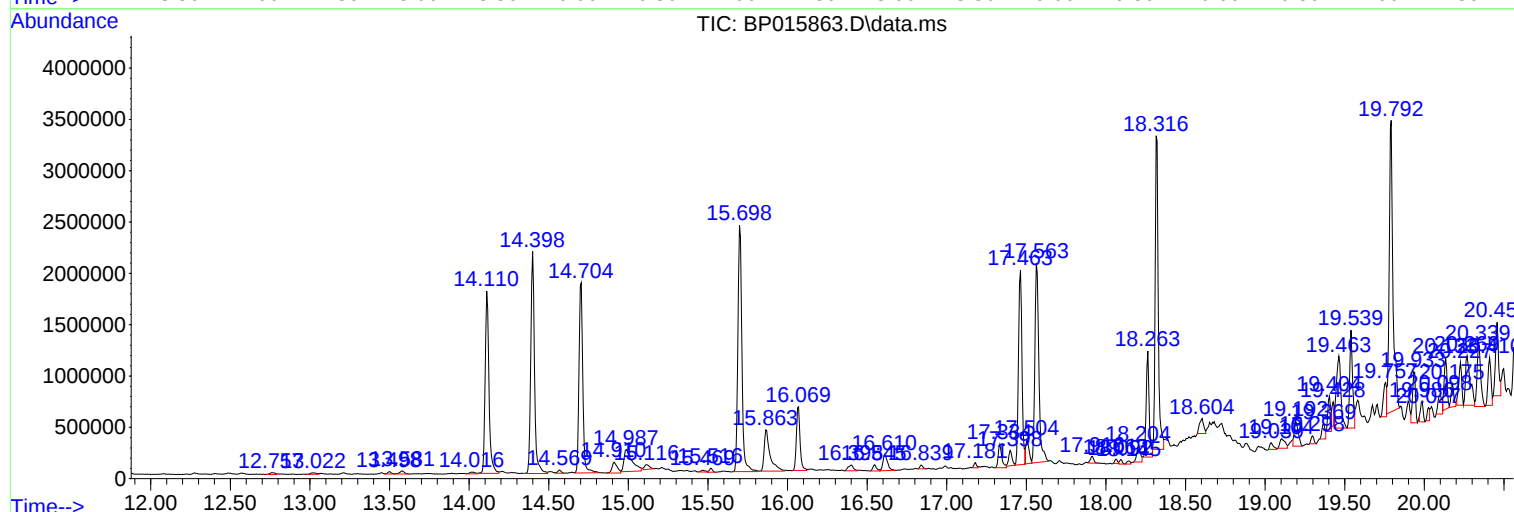
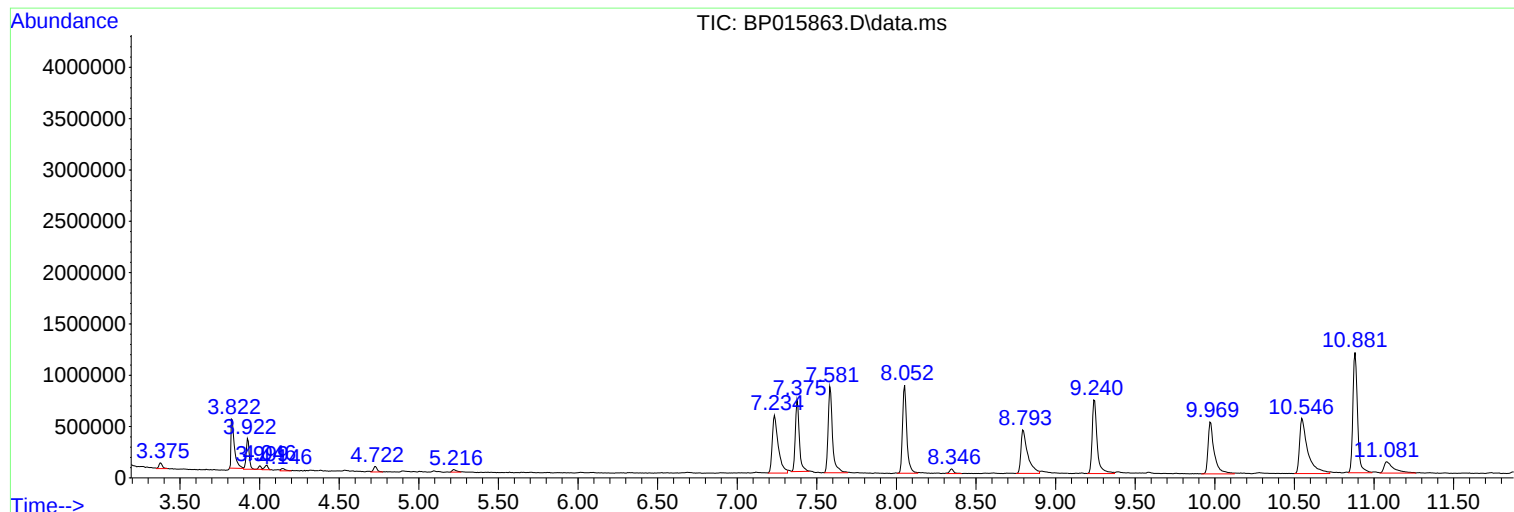
Sum of corrected areas: 119930089

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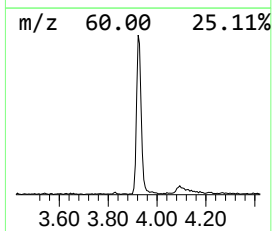
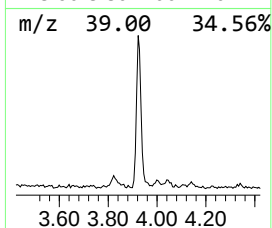
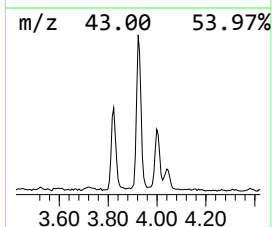
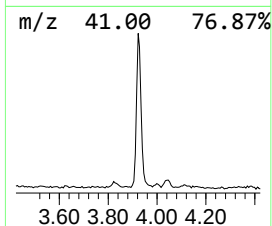
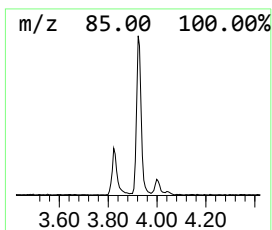
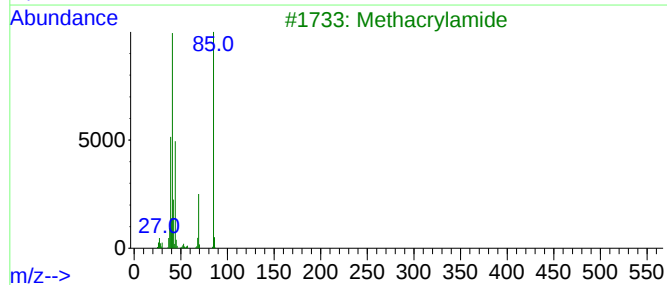
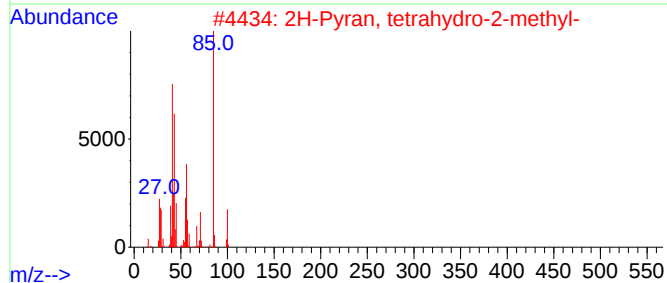
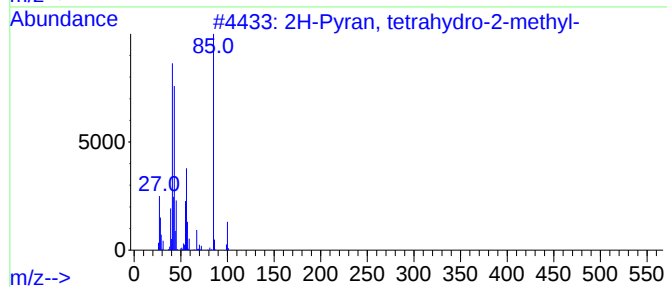
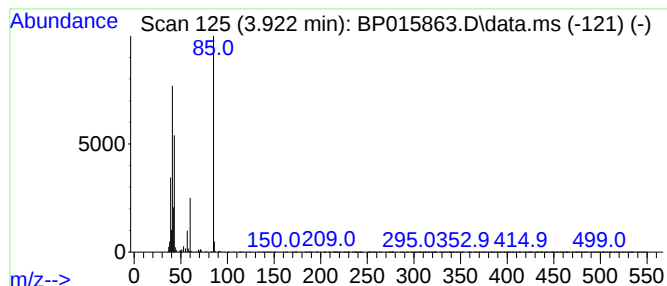
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.922	5.05 ng/ul	400583	1,4-Dichlorobenzene-d4	8.052

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	64
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	38
3	Methacrylamide			85	C4H7NO	000079-39-0	36
4	Oxetane, 2-ethyl-3-methyl-			100	C6H12O	053778-62-4	36
5	Methacrylamide			85	C4H7NO	000079-39-0	33



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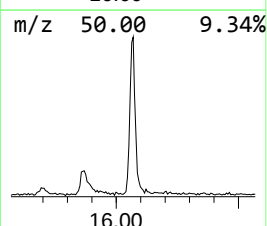
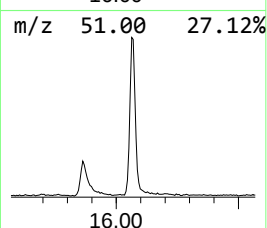
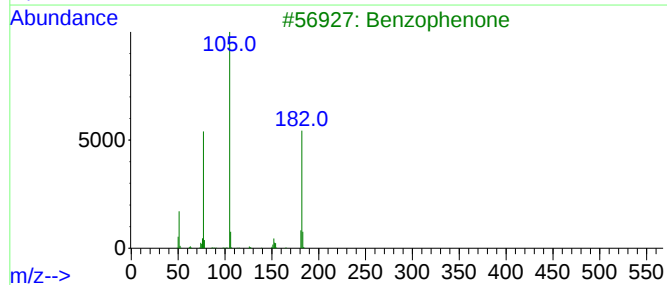
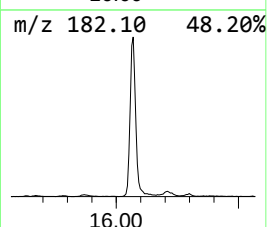
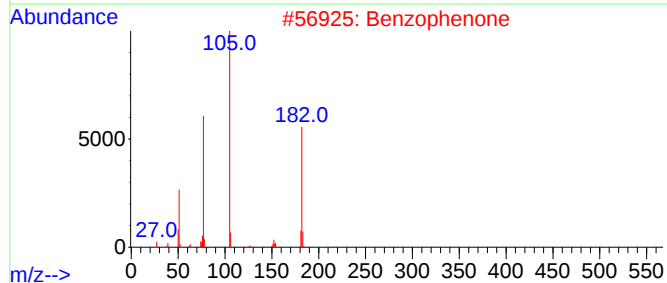
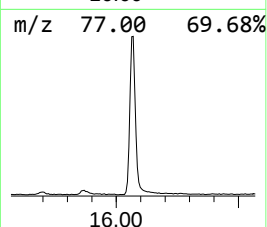
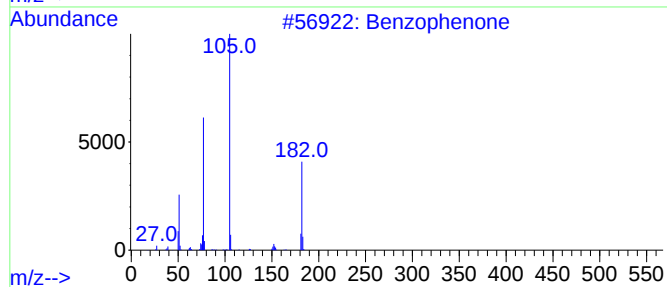
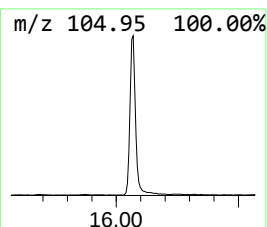
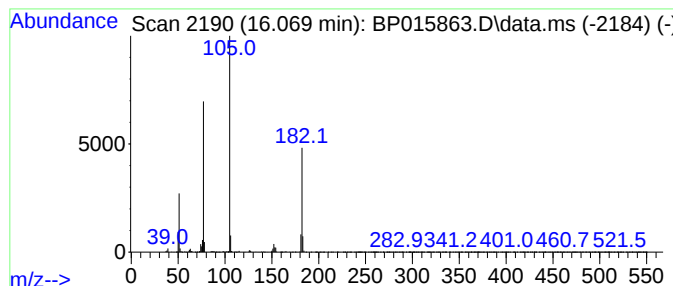
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	6.61 ng/ul	993609	Acenaphthene-d10	14.704

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	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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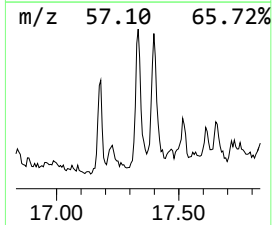
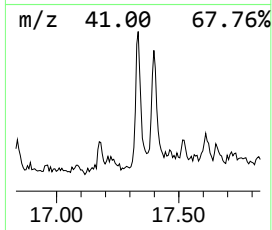
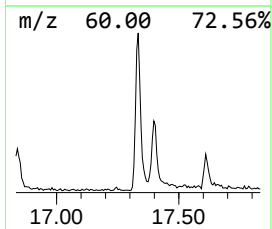
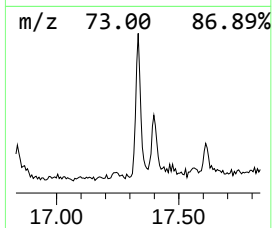
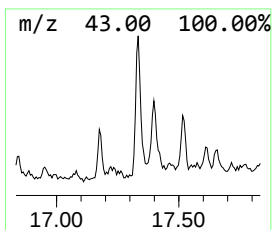
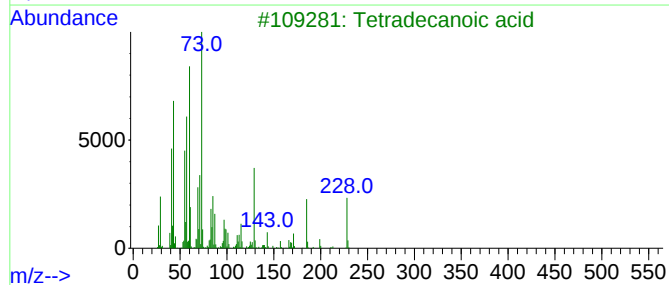
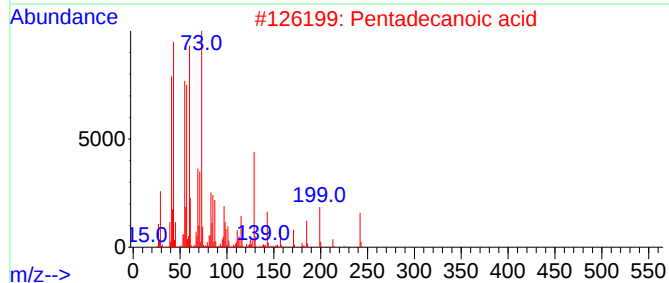
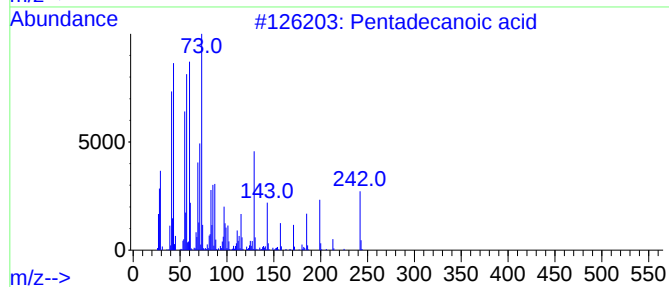
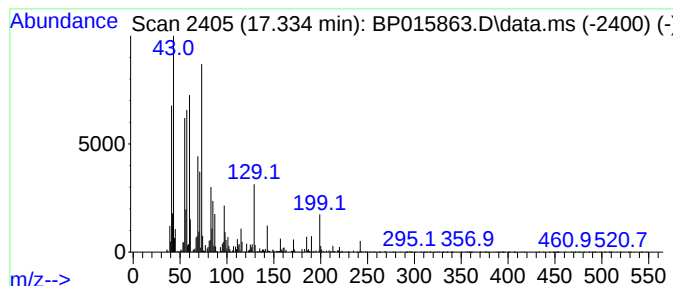
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Pentadecanoic acid Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.334	2.45 ng/ul	359535	Phenanthrene-d10	17.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecanoic acid	242	C15H30O2	001002-84-2	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	96
3	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	59
5	Heneicosanoic acid	326	C21H42O2	002363-71-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

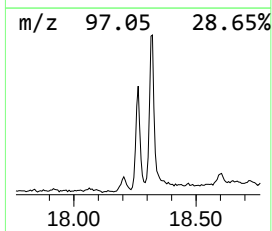
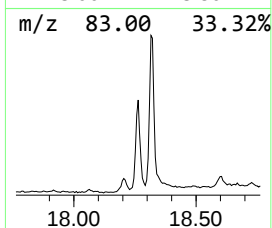
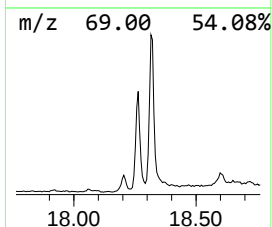
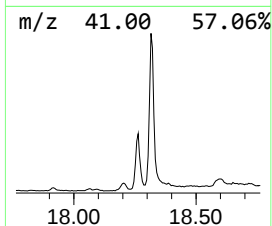
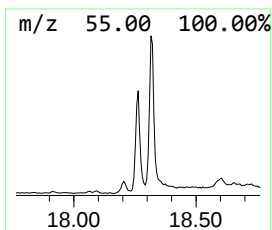
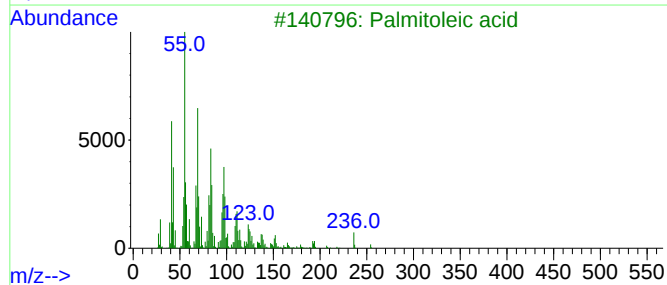
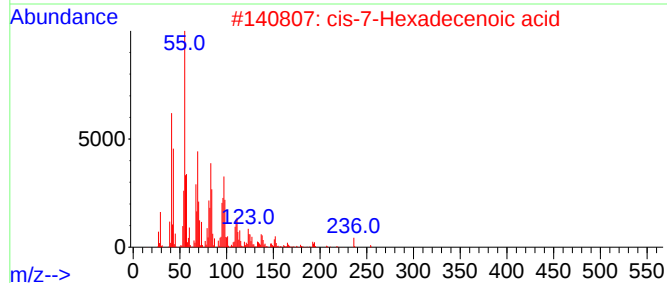
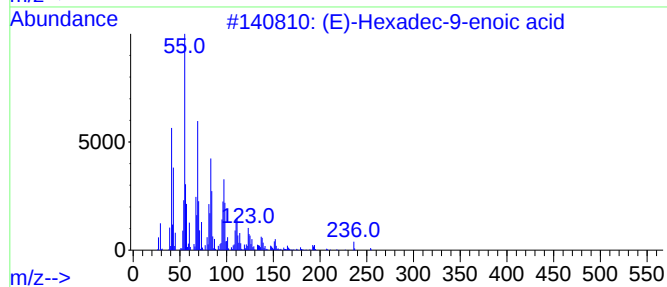
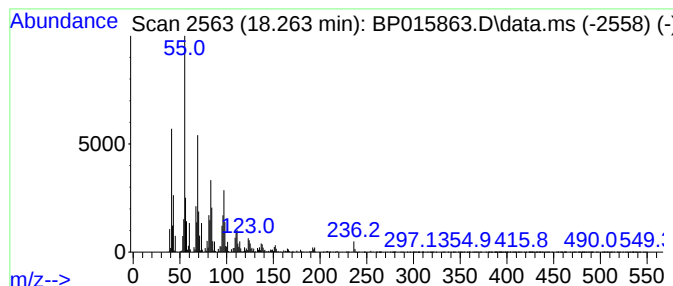
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (E)-Hexadec-9-enoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.263	8.98 ng/ul	1316120	Phenanthrene-d10	17.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
2	cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	99
3	Palmitoleic acid	254	C16H30O2	000373-49-9	94
4	9-Hexadecenoic acid	254	C16H30O2	002091-29-4	90
5	Myristoleic acid	226	C14H26O2	000544-64-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
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Instrument :
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ClientSampleId :
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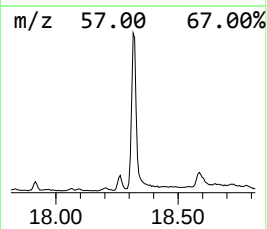
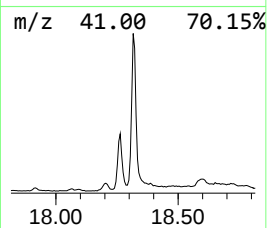
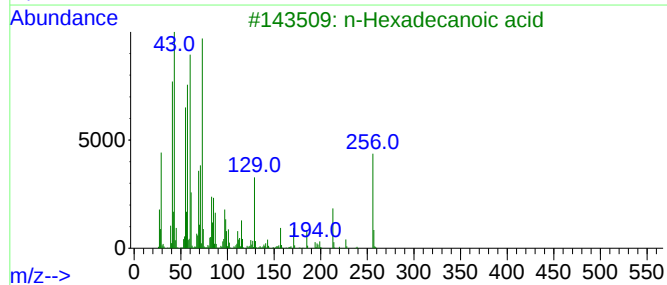
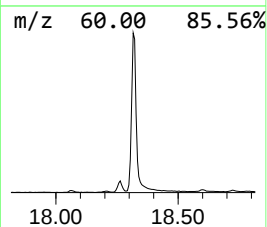
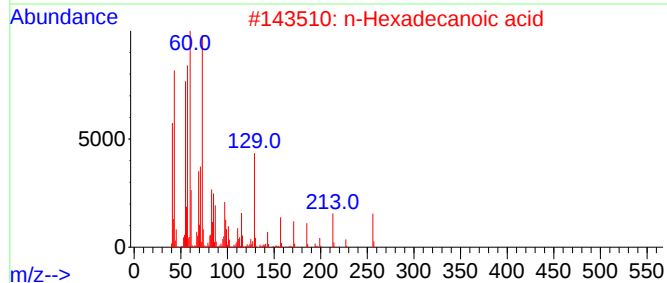
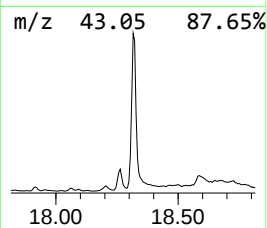
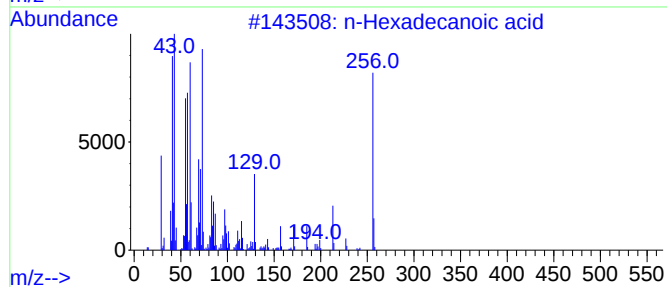
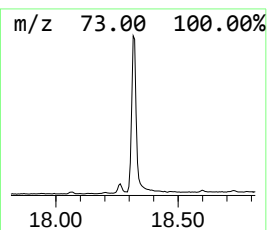
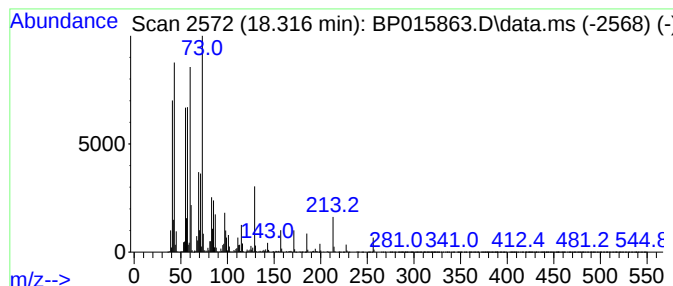
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	27.94 ng/ul	4093530	Phenanthrene-d10	17.463

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	96
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
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Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

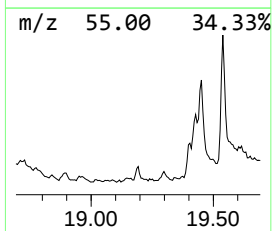
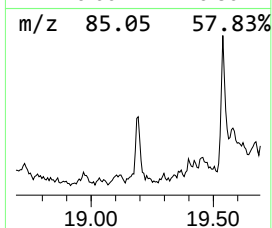
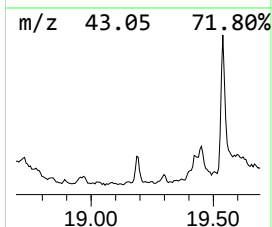
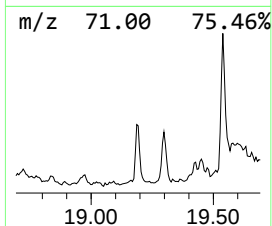
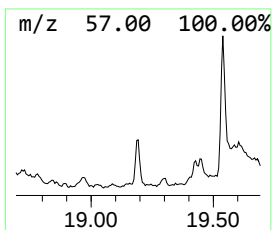
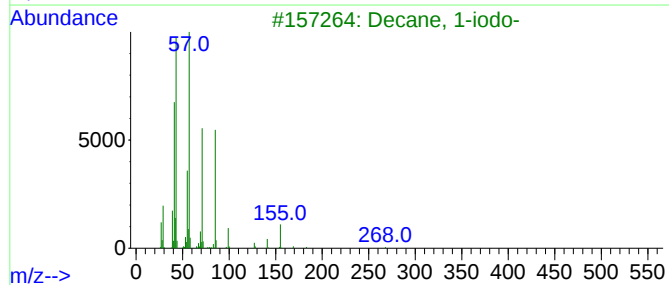
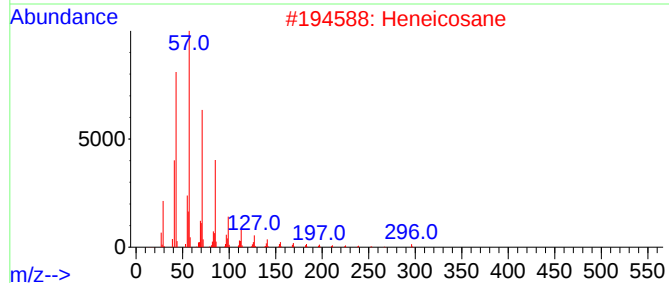
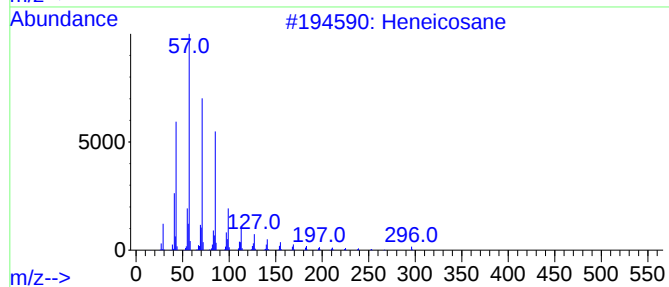
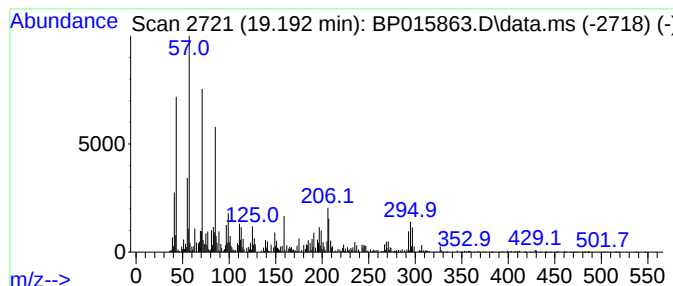
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.192	2.86 ng/ul	419137	Phenanthrene-d10		17.463	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	91
2	Heneicosane		296	C21H44	000629-94-7	64
3	Decane, 1-iodo-		268	C10H21I	002050-77-3	60
4	Octadecane		254	C18H38	000593-45-3	53
5	Tridecane		184	C13H28	000629-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
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Instrument :
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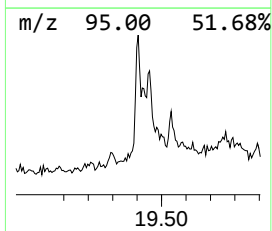
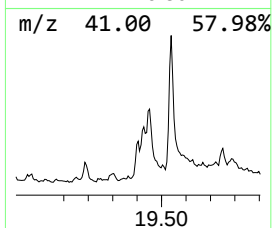
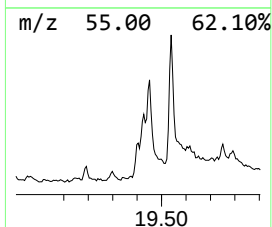
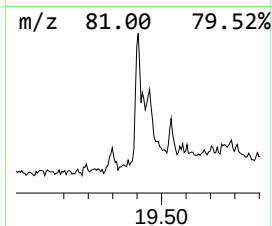
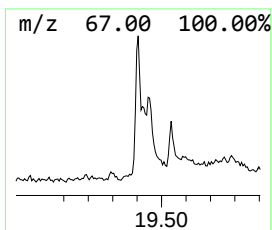
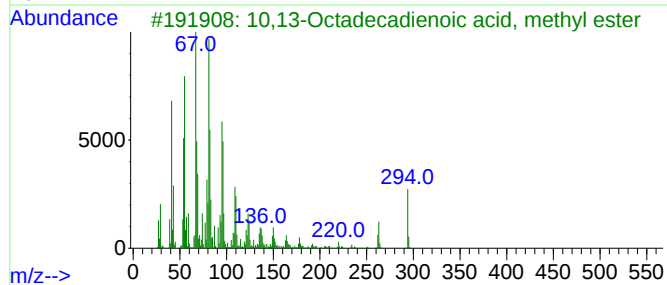
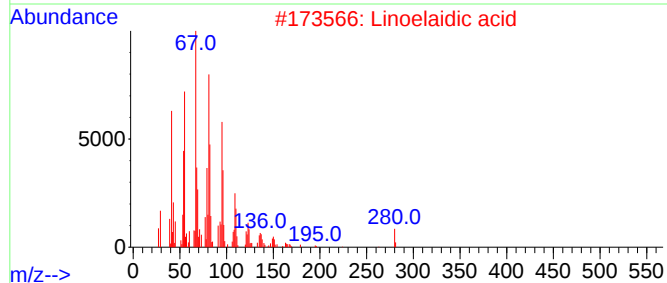
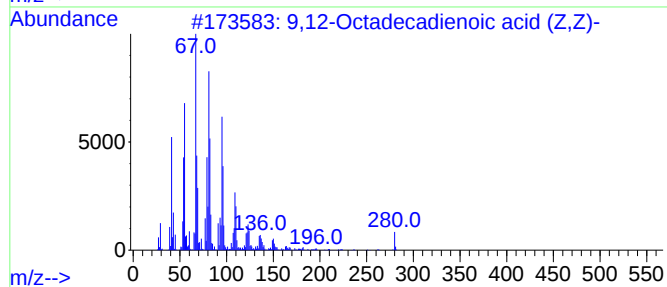
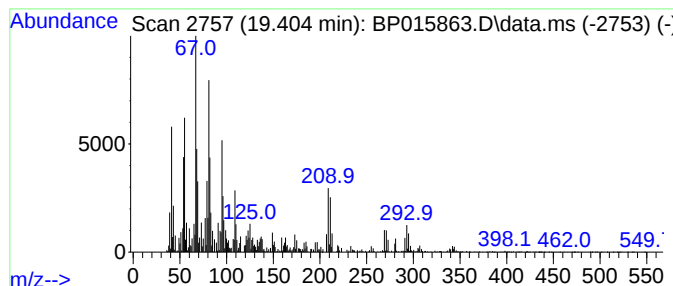
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 9,12-Octadecadienoic acid (... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.404	3.00 ng/ul	440229	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,12-Octadecadienoic acid (Z,Z)-	280	C18H32O2	000060-33-3	97
2			Linoelaidic acid	280	C18H32O2	000506-21-8	95
3			10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	92
4			7-Heptadecyne, 17-chloro-	270	C17H31Cl	056554-75-7	90
5			(9E,11E)-Octadecadienoic acid	280	C18H32O2	000544-71-8	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
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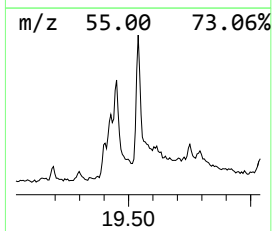
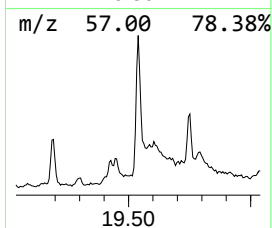
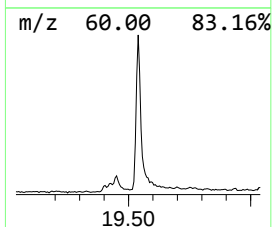
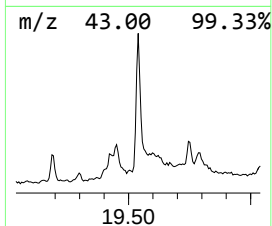
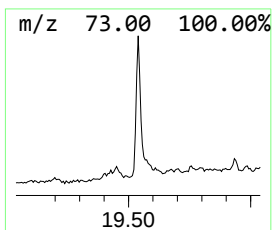
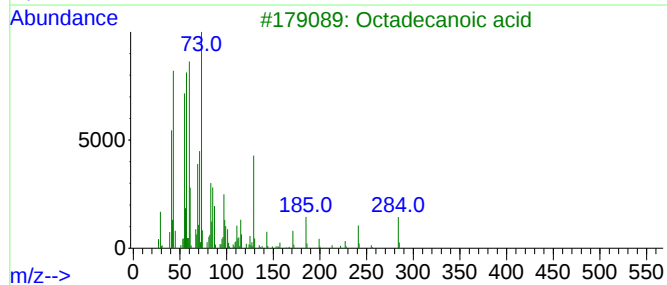
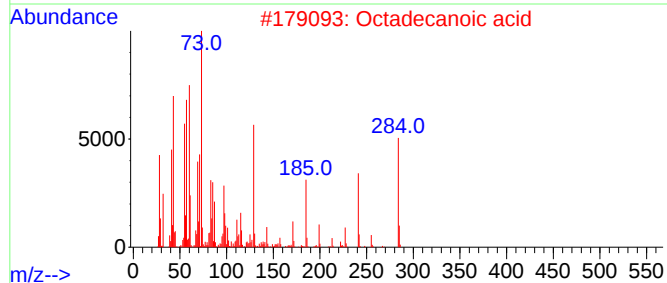
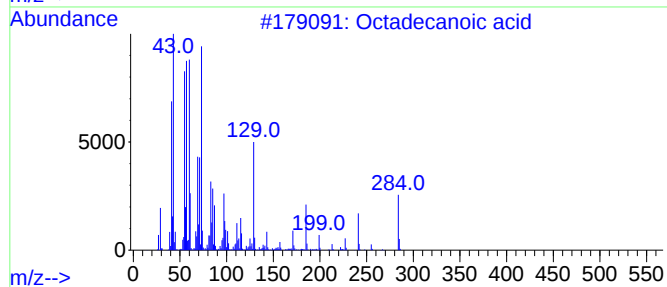
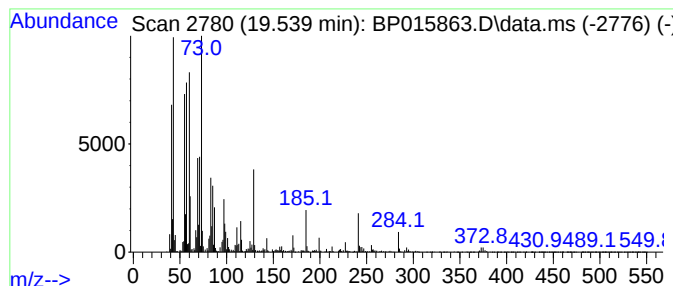
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	10.92 ng/ul	1229030	Chrysene-d12	21.551

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	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Misc :
ALS Vial : 29 Sample Multiplier: 1

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ClientSampleId :
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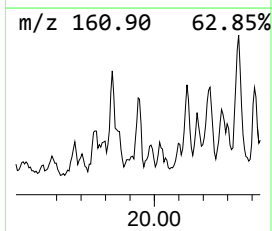
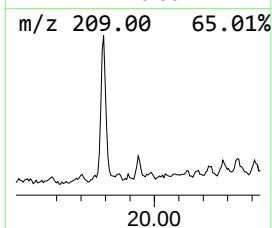
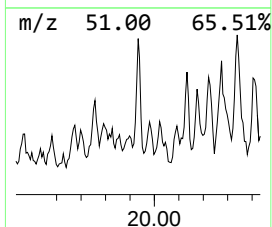
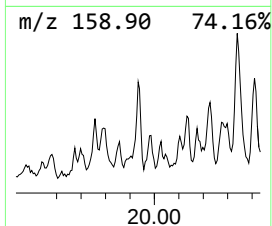
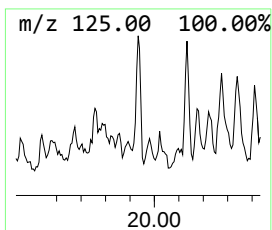
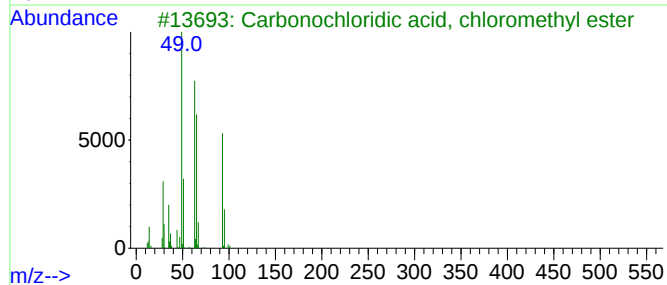
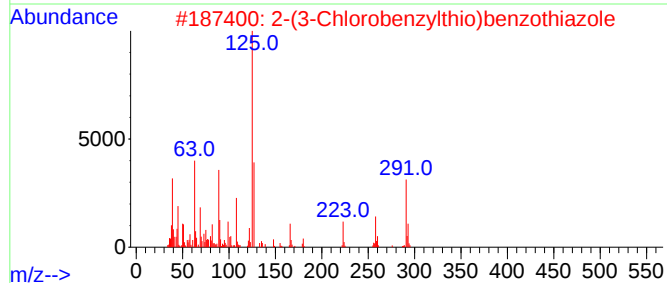
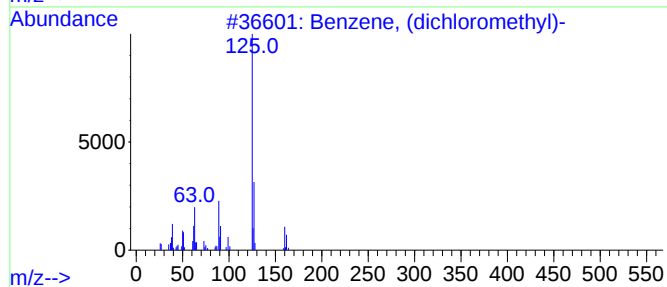
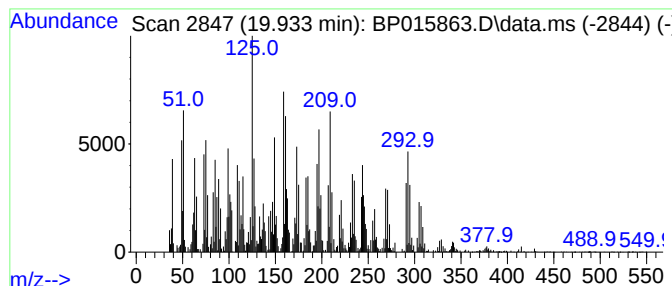
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-01 Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.933	5.50 ng/ul	618904	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (dichloromethyl)-	160	C7H6Cl2	000098-87-3	38
2			2-(3-Chlorobenzylthio)benzothiazole	291	C14H10ClNS2	342806-16-0	38
3			Carbonochloridic acid, chloromet...	128	C2H2Cl2O2	022128-62-7	32
4			2,6-Dichlorobenzylsulfonyl chloride	258	C7H5Cl3O2S	085952-31-4	25
5			5,5,6-Exo-8,9,10-hexachlorocamphene	340	C10H10Cl6	165820-10-0	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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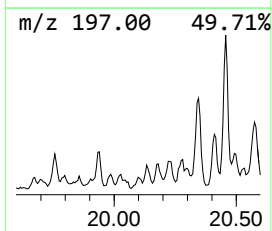
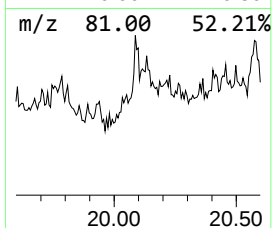
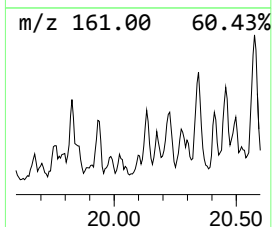
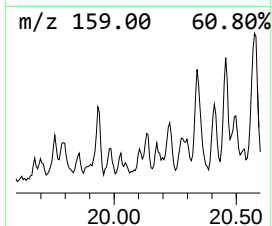
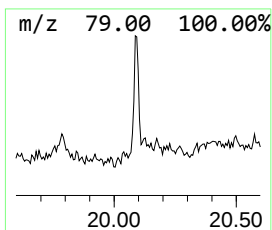
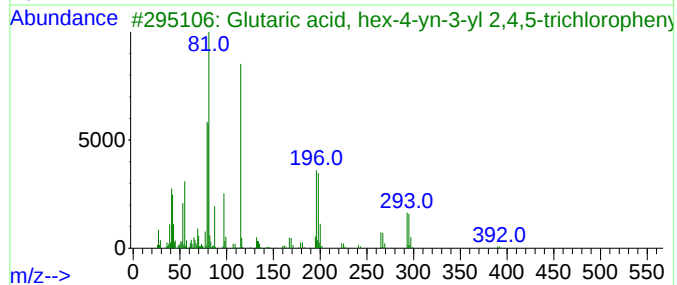
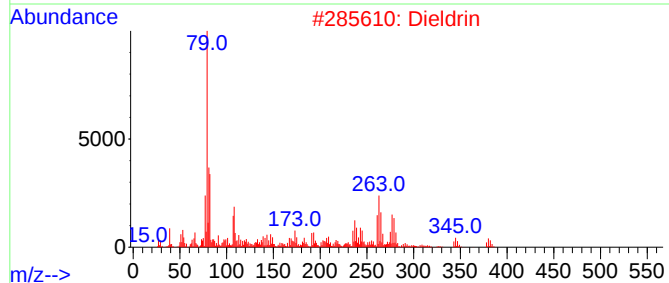
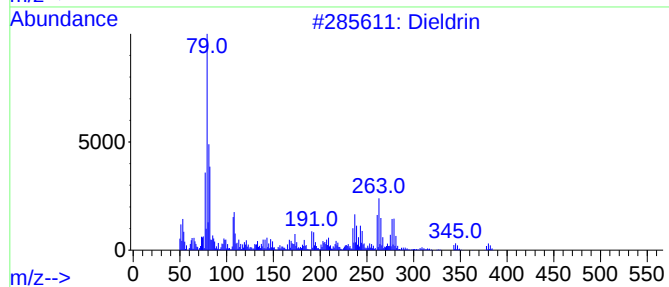
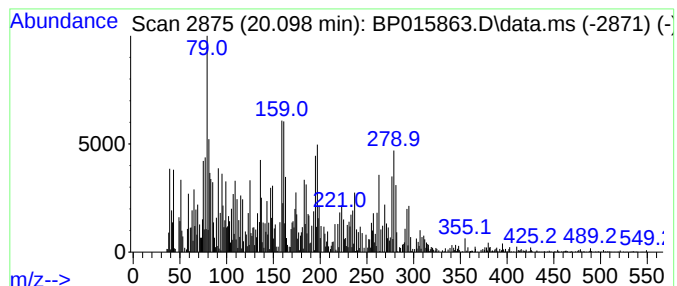
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-02 Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.098	3.02 ng/ul	339501	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dieldrin	378	C12H8Cl6O	000060-57-1	42
2			Dieldrin	378	C12H8Cl6O	000060-57-1	25
3			Glutaric acid, hex-4-yn-3-yl 2,4...	390	C17H17Cl3O4	1000392-16-1	10
4			Dieldrin	378	C12H8Cl6O	000060-57-1	9
5			S-(2,5-Dichloro-6-cyano-3-[(eth...	412	C12H10Cl2N2O2S4	194228-14-3	7



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFW8

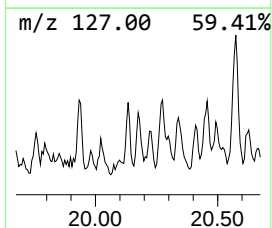
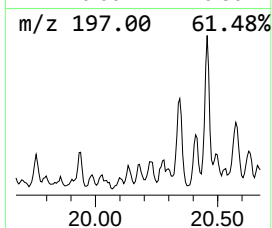
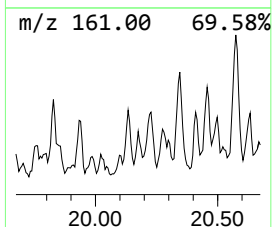
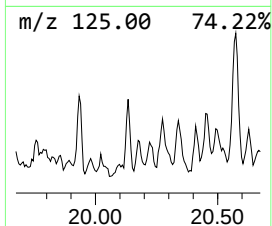
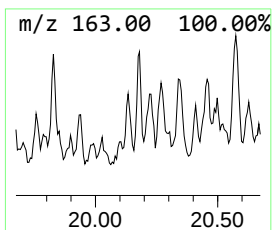
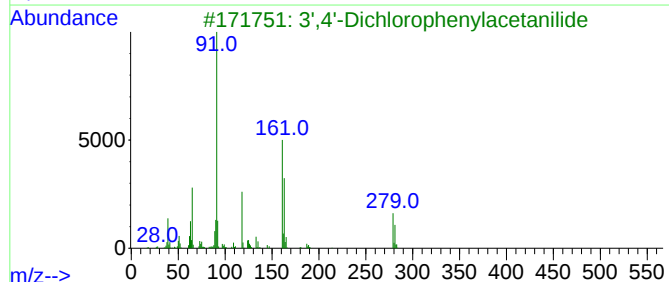
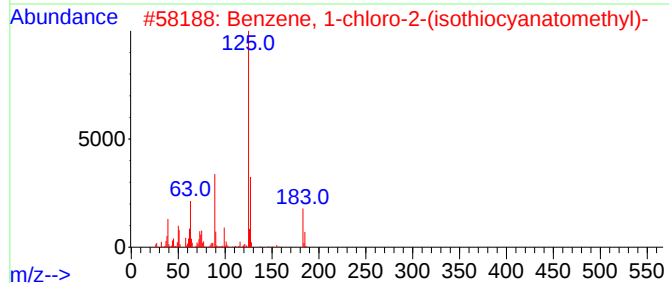
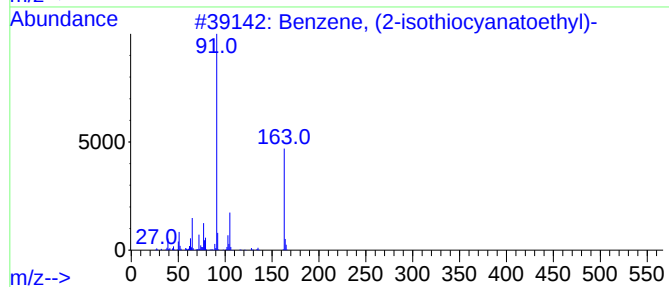
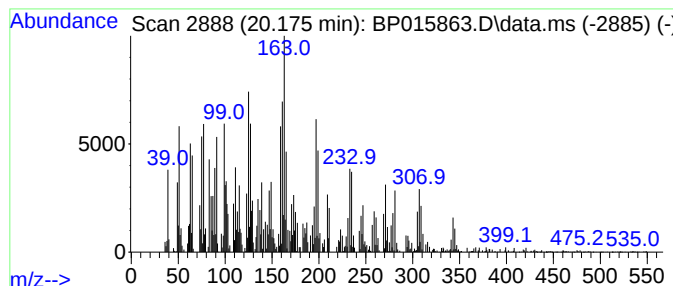
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-04 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.174	2.44 ng/ul	274259	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-isothiocyanatoethyl)-	163	C9H9NS	002257-09-2	22
2			Benzene, 1-chloro-2-(isothiocyan...	183	C8H6ClNS	018967-44-7	22
3			3',4'-Dichlorophenylacetanilide	279	C14H11Cl2NO	027816-82-6	18
4			2-Benzyloxy-5-bromo-4-(N,N-dimet...	307	C13H14BrN3O	205672-18-0	14
5			Succinic acid, 5-fluoro-2-nitrop...	319	C12H11ClFN06	1000357-98-0	12



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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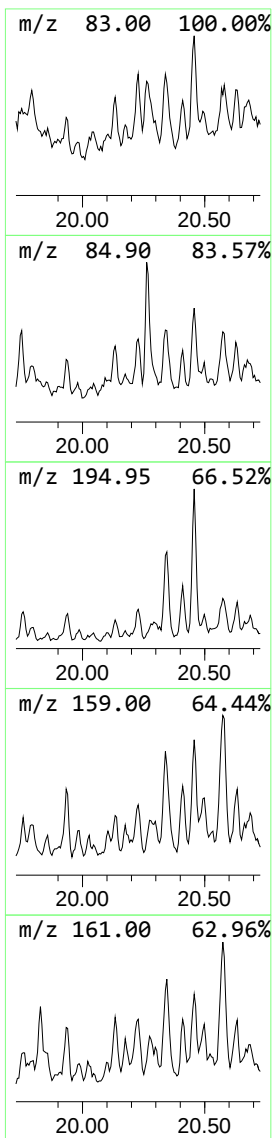
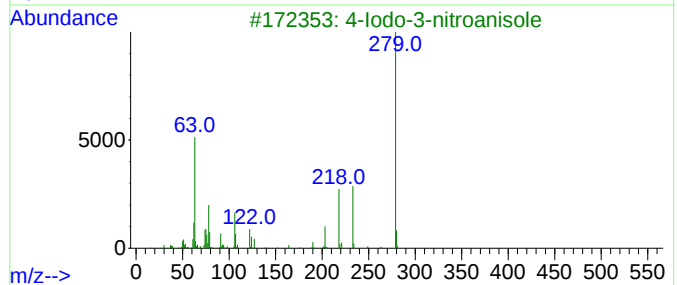
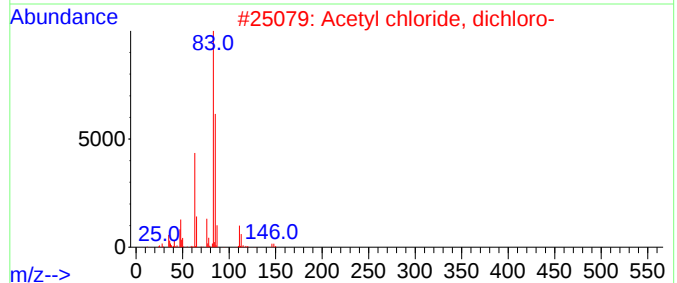
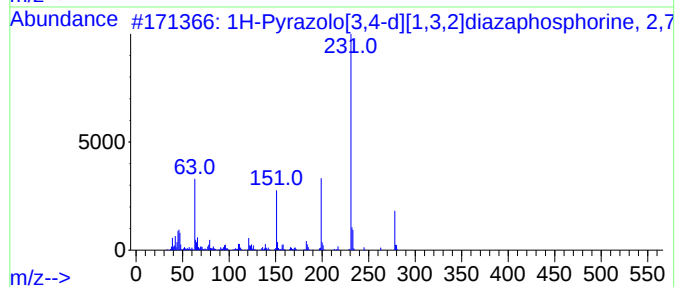
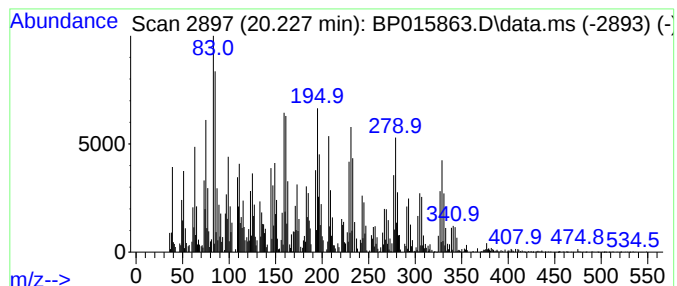
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-05 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.227	4.76 ng/ul	535570	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrazolo[3,4-d][1,3,2]diazaph...	278	C7H11N4PS3	1000350-06-9	35
2			Acetyl chloride, dichloro-	146	C2HCl3O	000079-36-7	25
3			4-Iodo-3-nitroanisole	279	C7H6INO3	058755-70-7	16
4			Phenol, 3-chloro-5-nitro-	173	C6H4ClNO3	000618-63-3	14
5			(1-Chloroethyl)trichlorosilane	196	C2H4Cl4Si	007787-82-8	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFW8

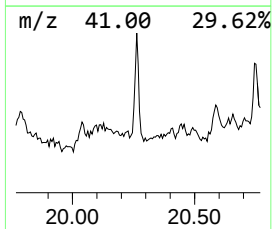
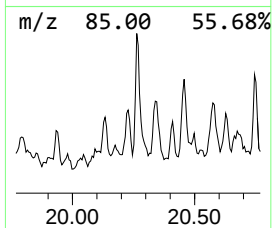
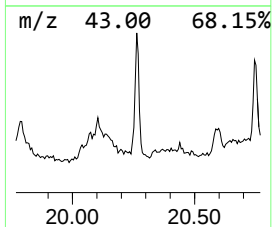
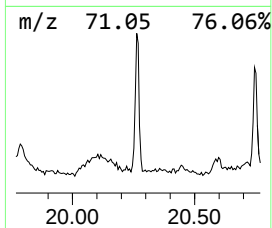
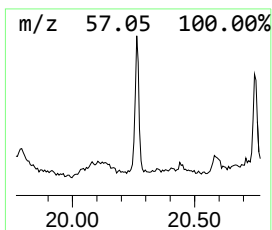
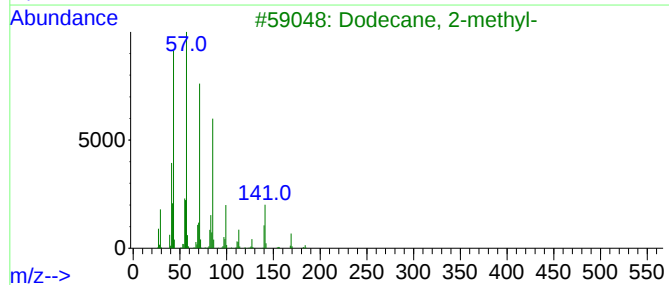
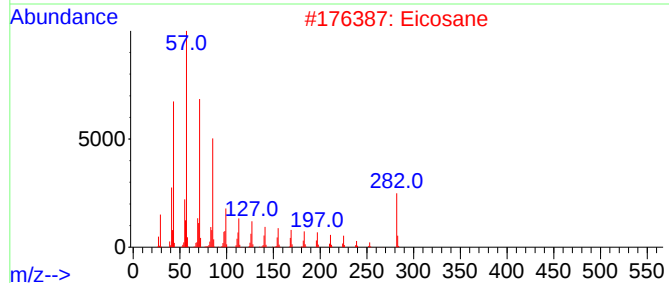
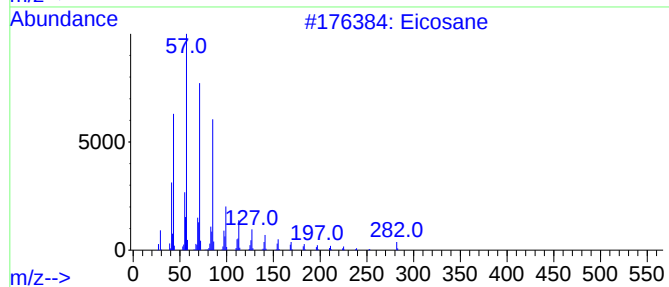
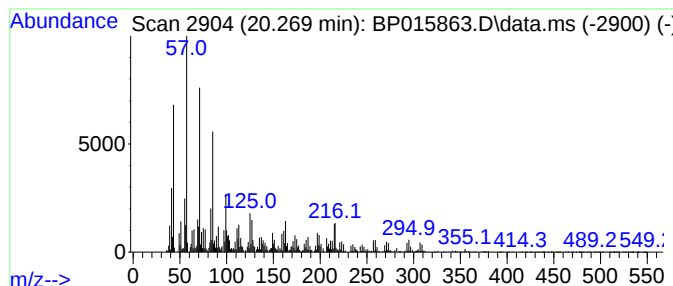
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.269	8.85 ng/ul	996115	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	86
2			Eicosane	282	C20H42	000112-95-8	76
3			Dodecane, 2-methyl-	184	C13H28	001560-97-0	76
4			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	70
5			Heptacosane	380	C27H56	000593-49-7	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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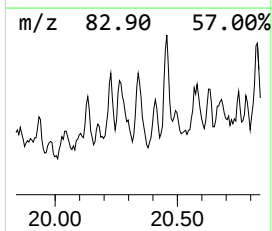
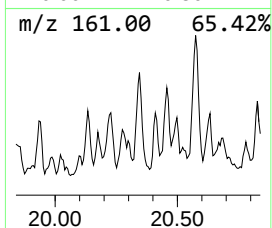
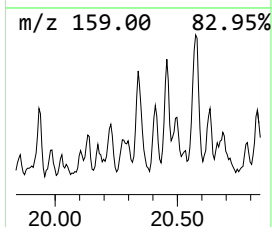
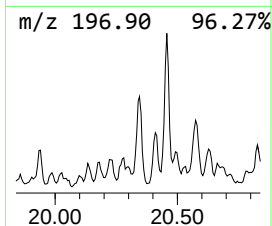
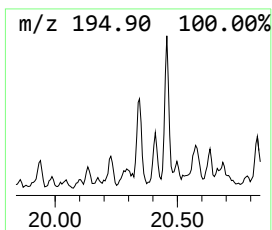
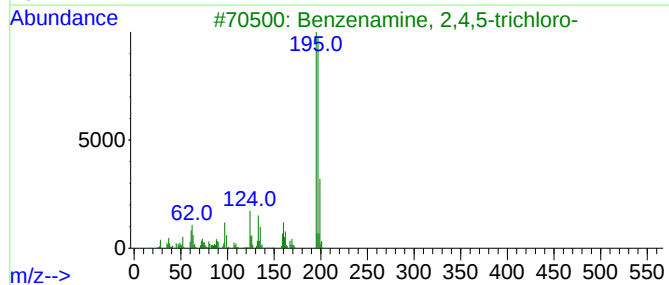
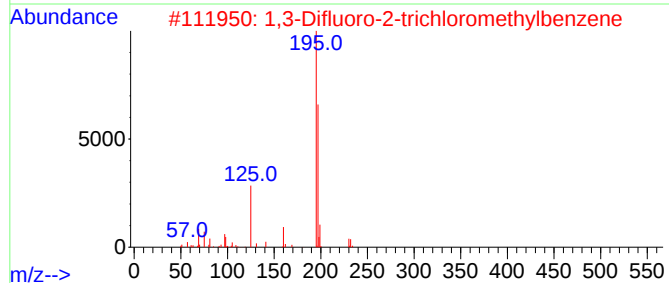
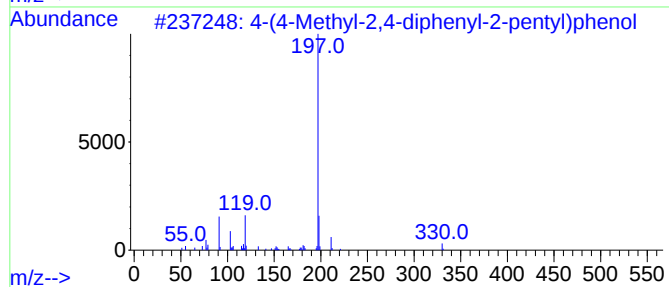
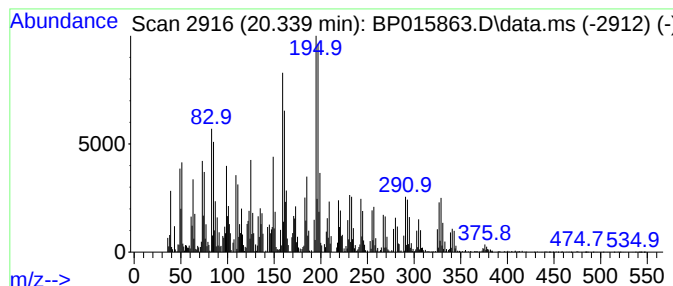
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-06 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.339	8.67 ng/ul	976217	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(4-Methyl-2,4-diphenyl-2-penty...	330	C24H26O	052379-26-7	44		
2	1,3-Difluoro-2-trichloromethylbe...	230	C7H3Cl3F2	093632-85-0	41		
3	Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	35		
4	Benzenamine, 2,3,4-trichloro-	195	C6H4Cl3N	000634-67-3	35		
5	4-Bromoindole	195	C8H6BrN	1000484-38-4	35		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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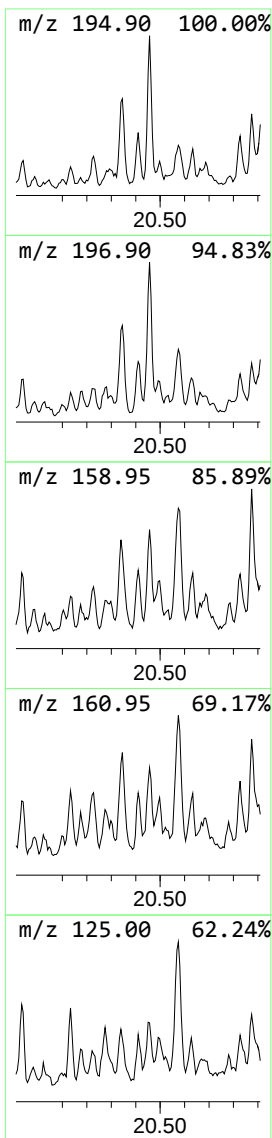
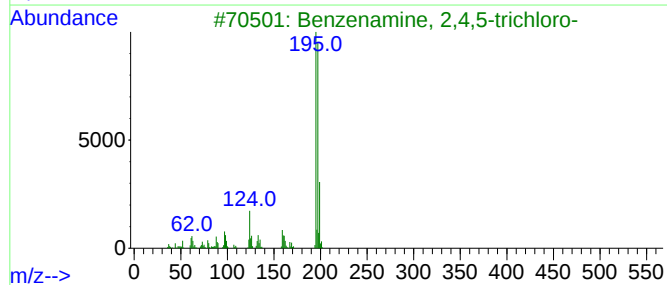
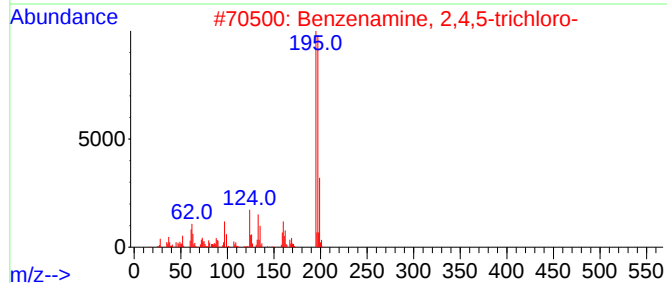
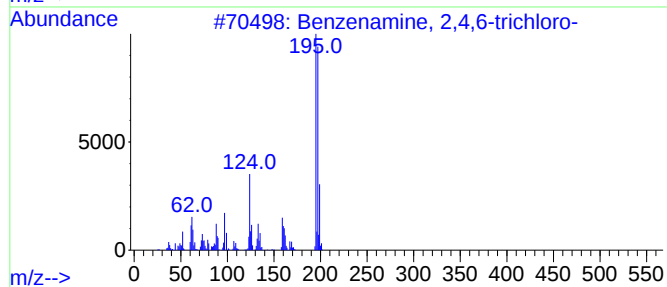
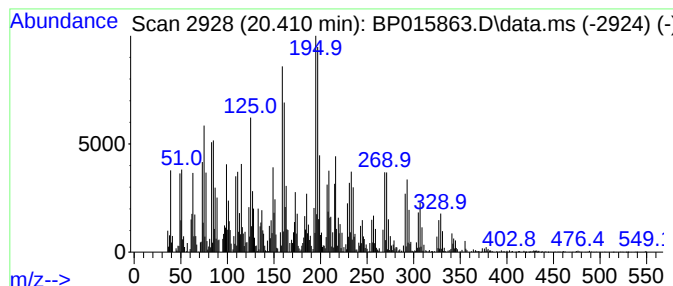
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 unknown-07 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.410	5.41 ng/ul	609522	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,4,6-trichloro-	195	C6H4Cl3N	000634-93-5	38
2			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	38
3			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	38
4			Benzenamine, 3,4,5-trichloro-	195	C6H4Cl3N	000634-91-3	38
5			4-Bromoindole	195	C8H6BrN	1000484-38-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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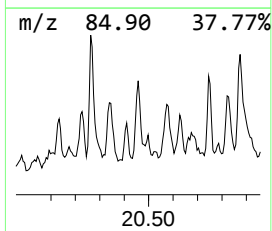
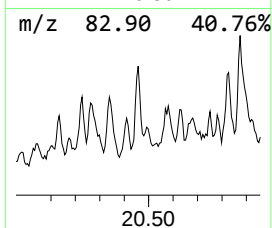
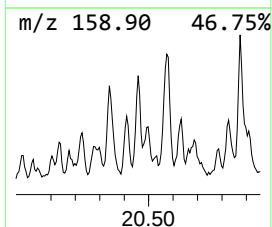
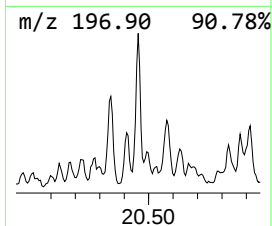
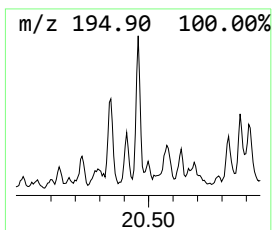
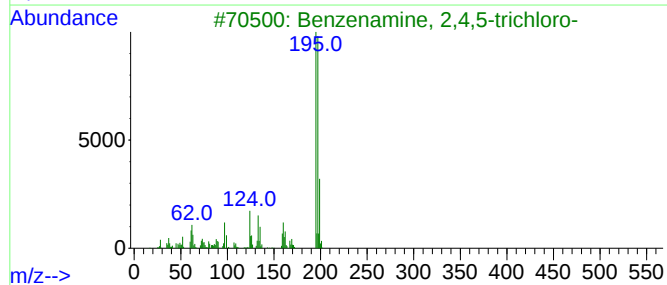
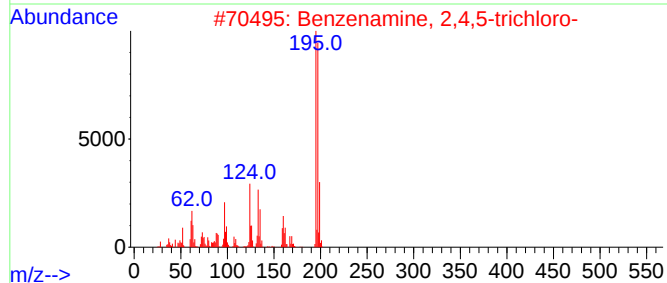
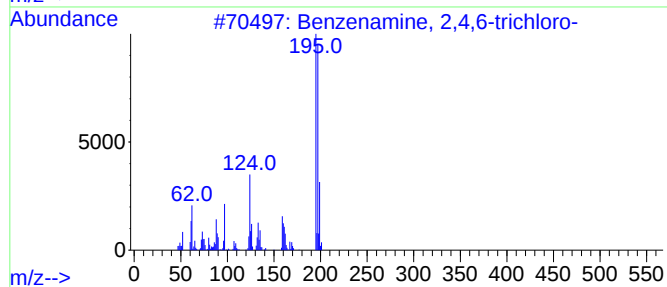
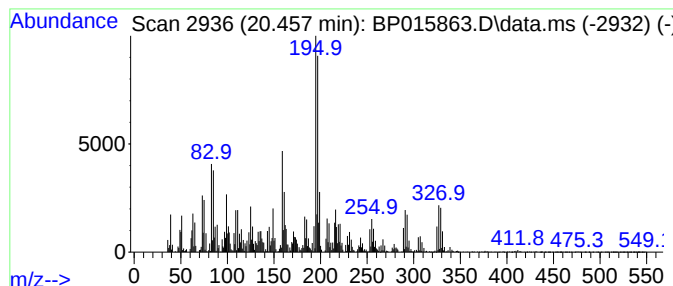
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Benzenamine, 2,4,6-trichloro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.457	8.74 ng/ul	984361	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,4,6-trichloro-	195	C6H4Cl3N	000634-93-5	60
2			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	55
3			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	55
4			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	55
5			Benzenamine, 2,3,4-trichloro-	195	C6H4Cl3N	000634-67-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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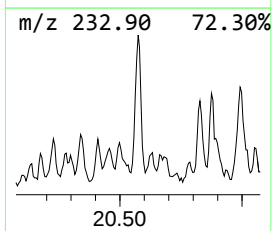
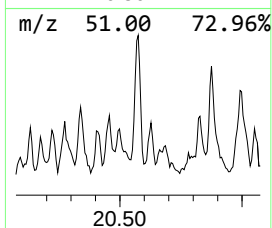
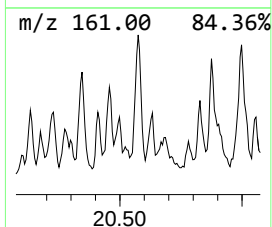
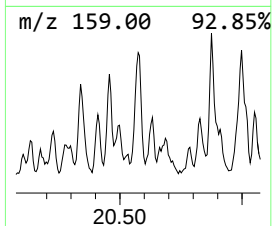
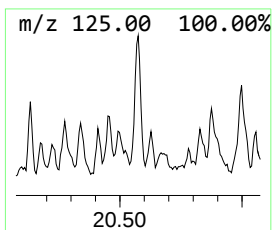
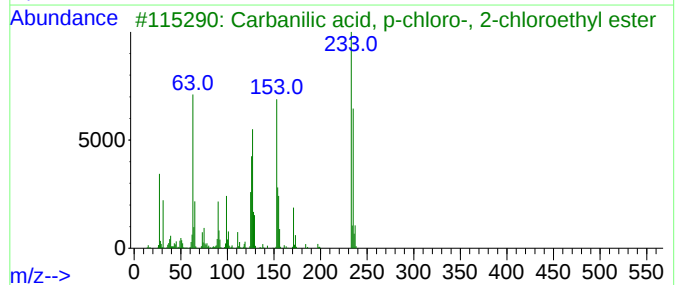
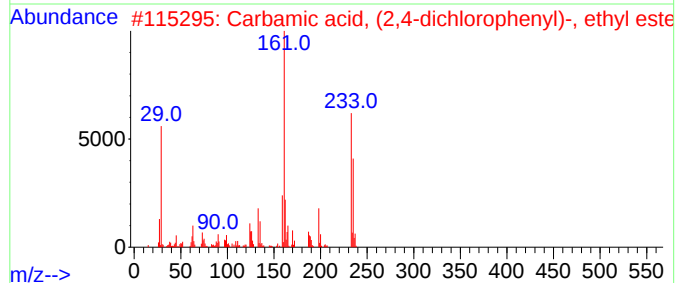
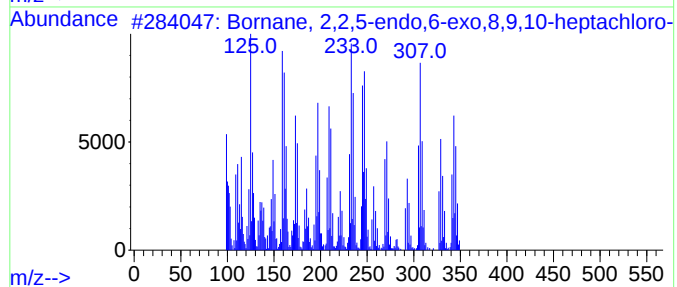
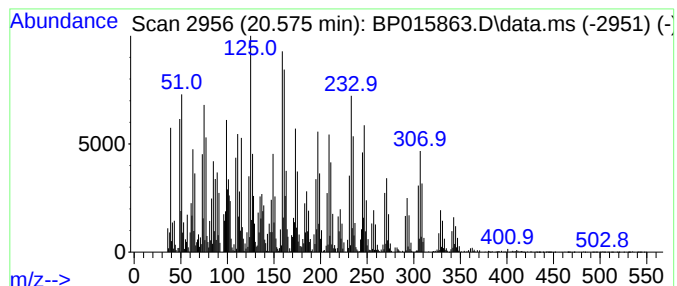
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Bornane, 2,2,5-endo,6-exo,8... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.575	15.03 ng/ul	1691400	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bornane, 2,2,5-endo,6-exo,8,9,10...	376	C10H11Cl7	051775-36-1	76
2			Carbamic acid, (2,4-dichlorophen...	233	C9H9Cl2NO2	006333-37-5	45
3			Carbanilic acid, p-chloro-, 2-ch...	233	C9H9Cl2NO2	025217-18-9	44
4			Carbamic acid, (3,4-dichlorophen...	233	C9H9Cl2NO2	007159-94-6	25
5			2-Chloro-5-nitrophenol	173	C6H4ClNO3	000619-10-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFw8

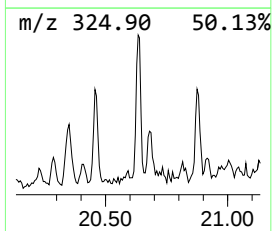
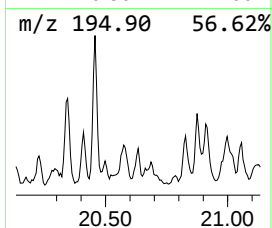
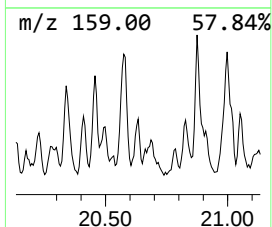
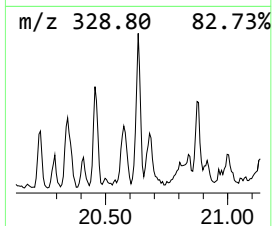
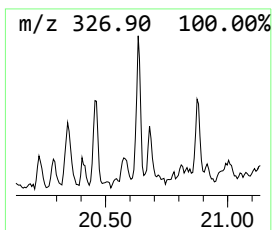
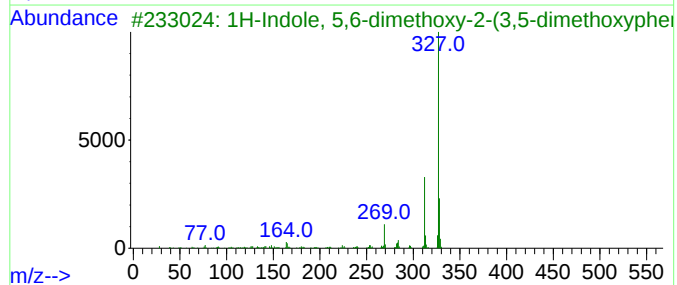
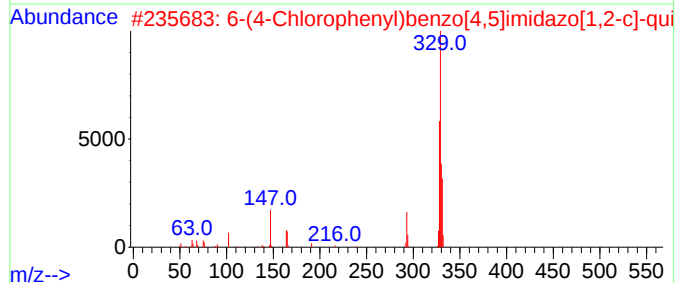
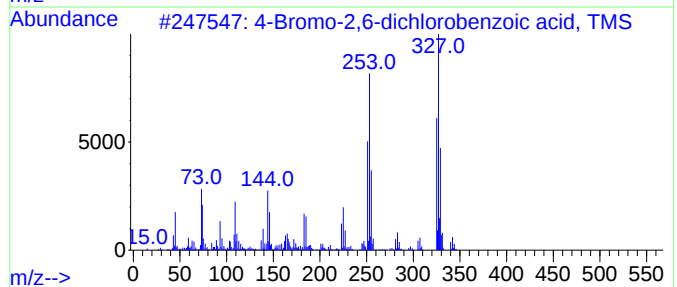
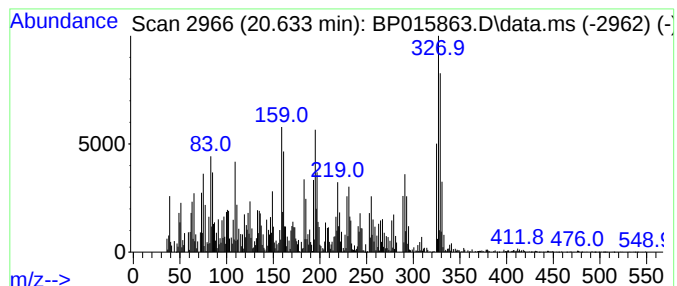
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown-08 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.633	5.03 ng/ul	566320	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Bromo-2,6-dichlorobenzoic acid...	340	C10H11BrCl2O2Si	1000497-68-4	22
2			6-(4-Chlorophenyl)benzo[4,5]imid...	329	C20H12ClN3	105997-06-6	14
3			1H-Indole, 5,6-dimethoxy-2-(3,5-...	327	C19H21NO4	156785-69-2	11
4			1,2-bis(4-Methoxyphenyl)indole	329	C22H19NO2	1262543-81-6	11
5			1H-Pyrazole-5-carboxamide, 3-cyc...	327	C20H29N3O	1000350-11-6	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFW8

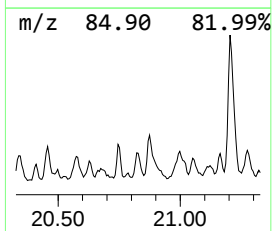
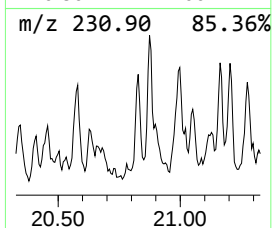
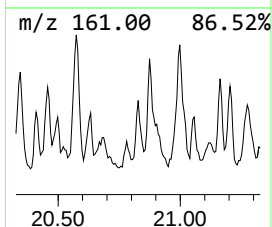
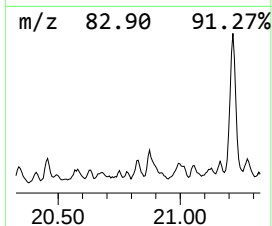
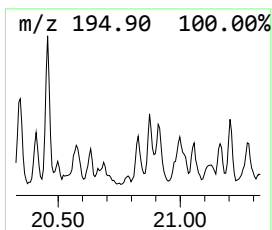
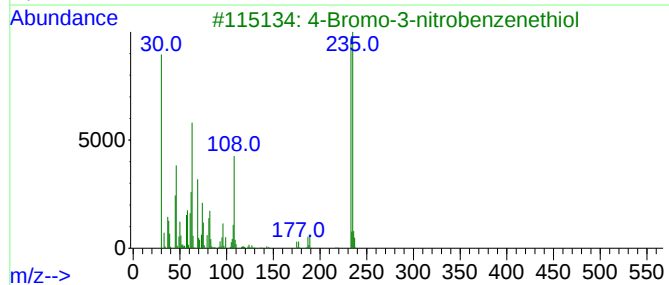
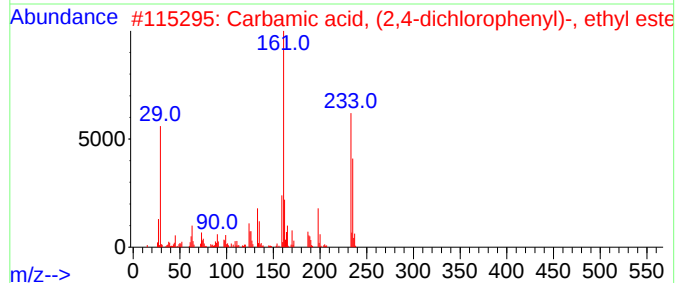
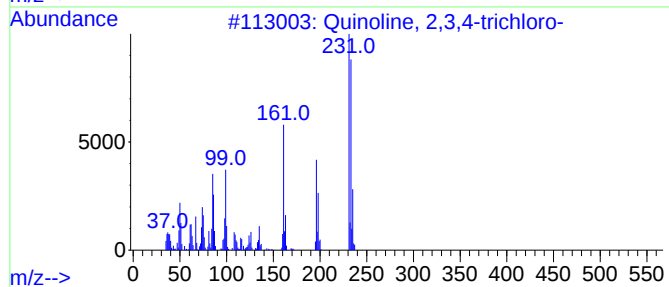
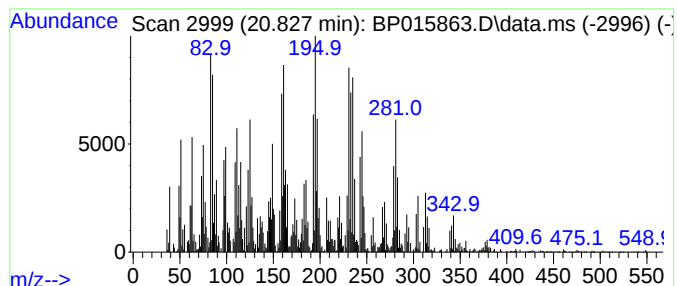
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 unknown-09 Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.827	4.73 ng/ul	532995	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Quinoline, 2,3,4-trichloro-	231	C9H4Cl3N	040335-02-2	42
2			Carbamic acid, (2,4-dichlorophen...	233	C9H9Cl2NO2	006333-37-5	35
3			4-Bromo-3-nitrobenzenethiol	233	C6H4BrNO2S	1000459-09-8	25
4			4-(4-Bromophenyl)pyridine	233	C11H8BrN	039795-60-3	25
5			Carbamic acid, (3,4-dichlorophen...	233	C9H9Cl2NO2	007159-94-6	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFW8

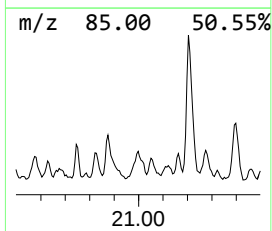
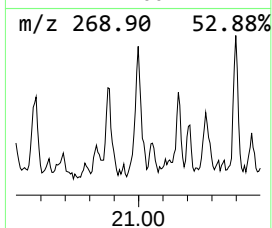
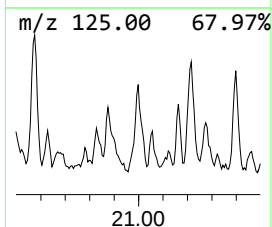
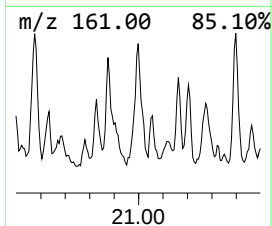
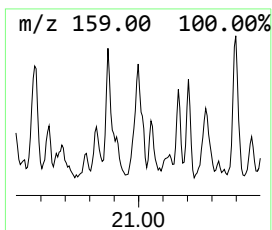
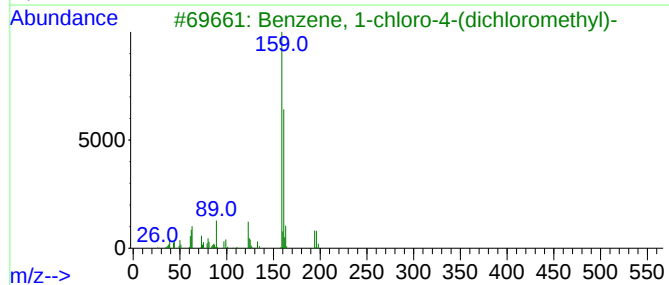
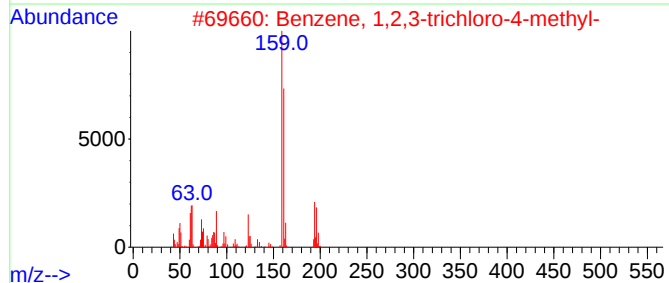
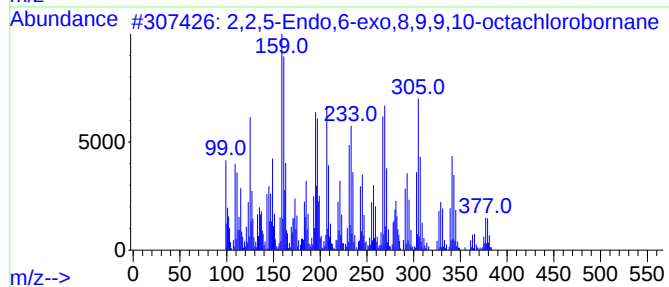
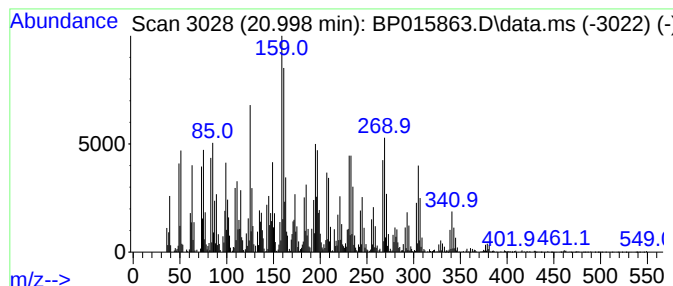
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 2,2,5-Endo,6-exo,8,9,9,10-o... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.998	13.66 ng/ul	1537420	Chrysene-d12	21.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,5-Endo,6-exo,8,9,9,10-octach...	410	C10H10Cl8	058002-19-0	90	
2	Benzene, 1,2,3-trichloro-4-methyl-	194	C7H5Cl3	007359-72-0	41	
3	Benzene, 1-chloro-4-(dichloromet...	194	C7H5Cl3	013940-94-8	25	
4	Bis-2-chloroethyl-2-chloroethylp...	268	C6H12Cl3O3P	006294-34-4	18	
5	Dimethylgermanium dichloride	174	C2H6Cl2Ge	001529-48-2	14	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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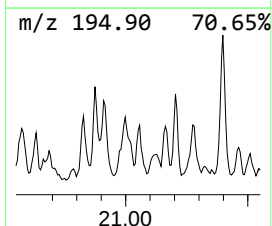
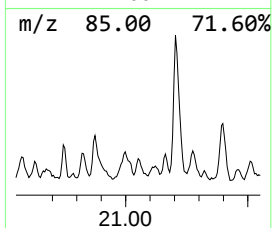
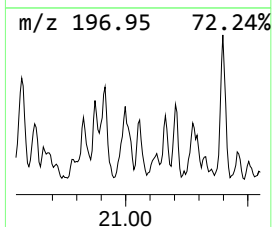
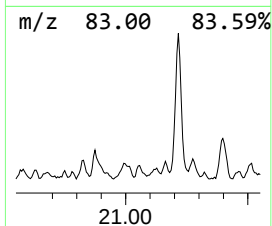
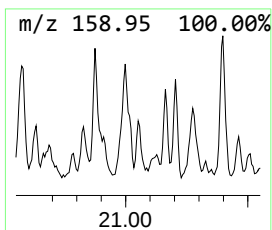
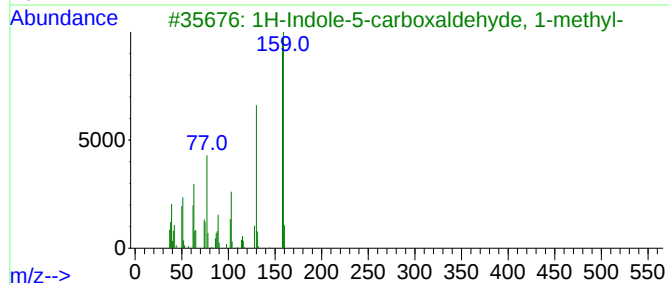
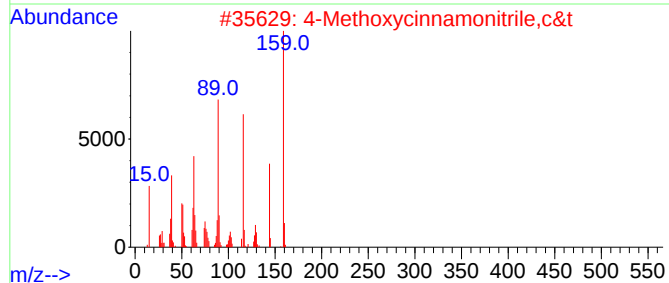
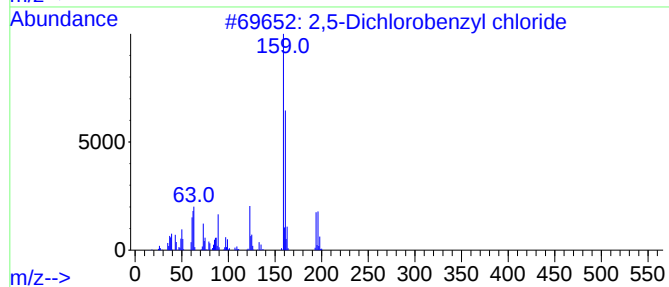
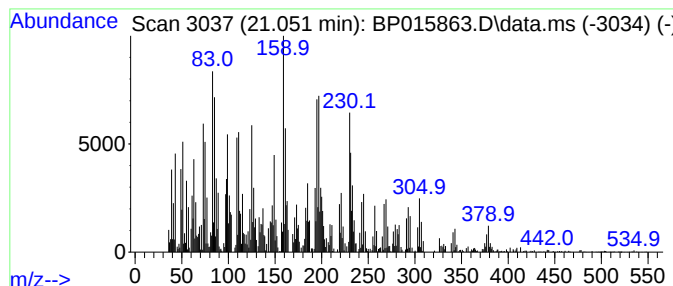
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-11 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.051	4.47 ng/ul	503431	Chrysene-d12	21.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,5-Dichlorobenzyl chloride	194	C7H5Cl3	002745-49-5	46
2		4-Methoxycinnamitrile,c&t	159	C10H9NO	028446-68-6	30
3		1H-Indole-5-carboxaldehyde, 1-me...	159	C10H9NO	090923-75-4	27
4		Propane, 1,1,1,3-tetrachloro-	180	C3H4Cl4	001070-78-6	27
5		3-endo,5-endo-dichloro-7,7-bis-c...	340	C10H10Cl6	1000365-99-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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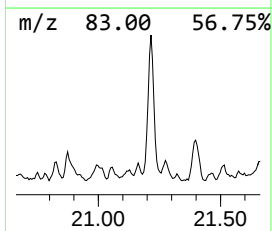
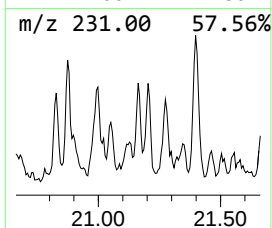
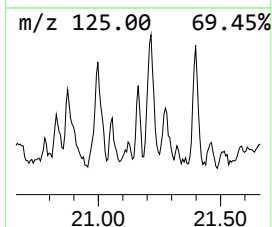
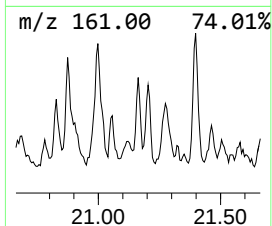
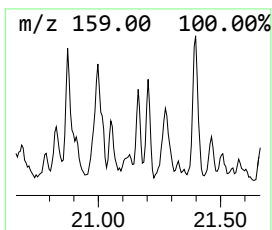
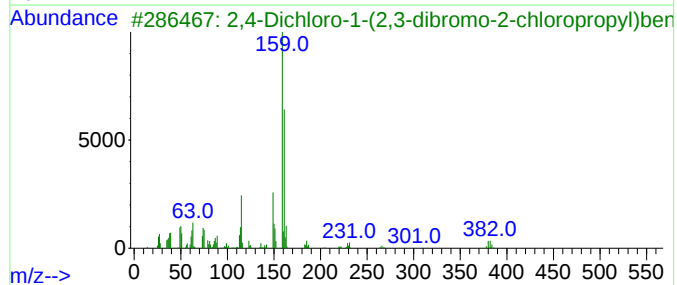
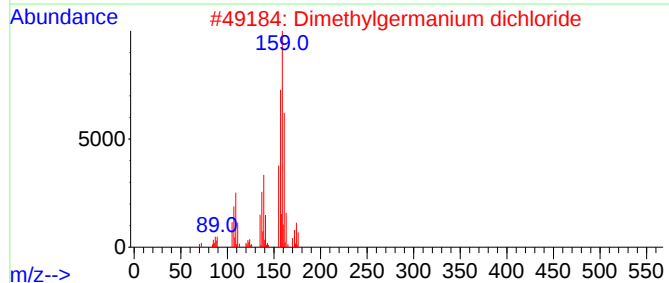
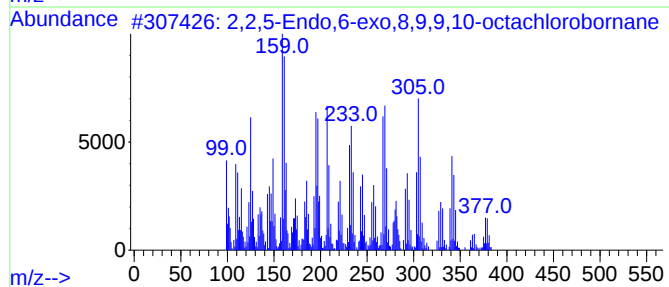
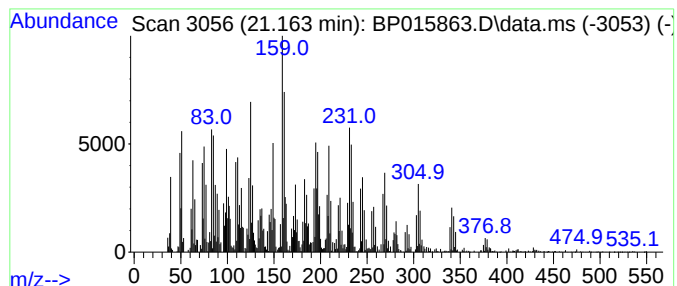
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 unknown-12 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.163	5.78 ng/ul	650145	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,5-Endo,6-exo,8,9,9,10-octach...	410	C10H10Cl8	058002-19-0	42		
2	Dimethylgermanium dichloride	174	C2H6Cl2Ge	001529-48-2	30		
3	2,4-Dichloro-1-(2,3-dibromo-2-ch...	378	C9H7Br2Cl3	1000444-78-6	15		
4	Benzene, 1,2,3-trichloro-4-methyl-	194	C7H5Cl3	007359-72-0	15		
5	Benzene, 1,3-dichloro-5-(chlorom...	194	C7H5Cl3	003290-06-0	15		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

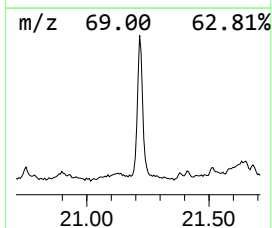
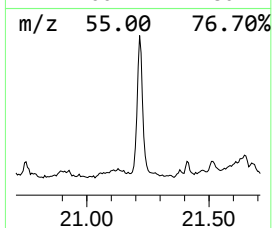
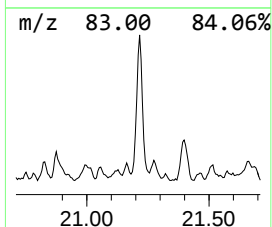
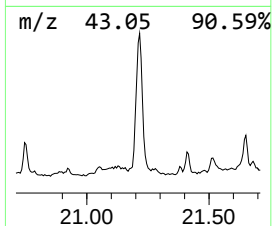
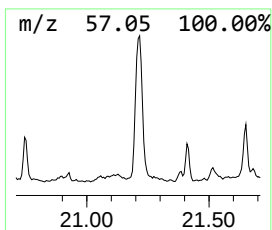
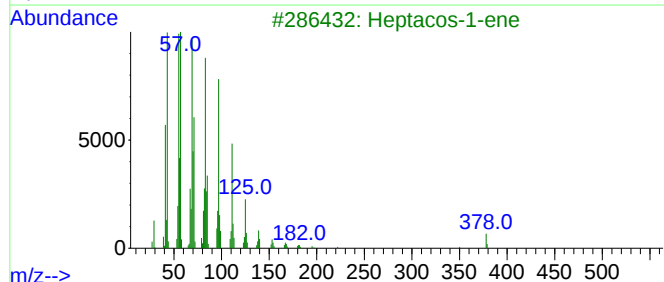
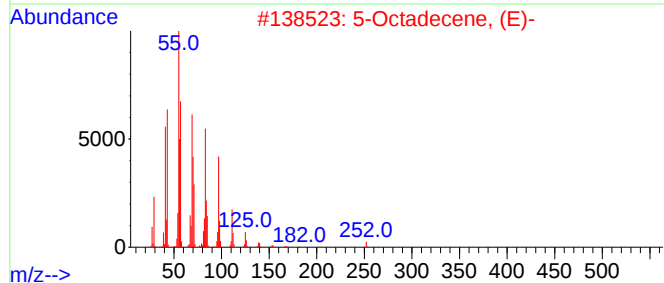
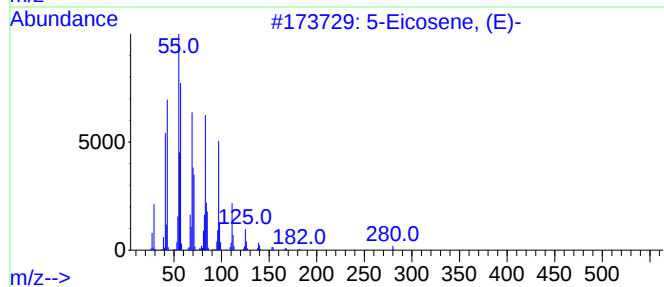
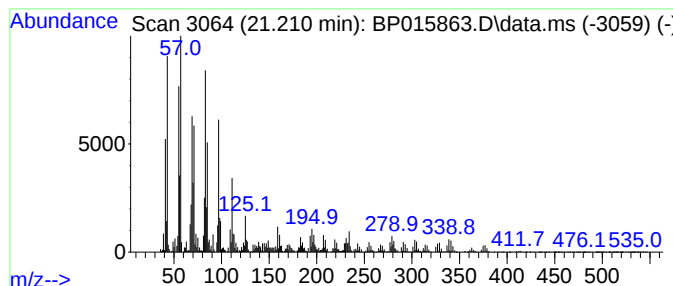
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.210	28.84 ng/ul	3246300	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-	280	C20H40	074685-30-6	93
2	5-Octadecene, (E)-	252	C18H36	007206-21-5	89
3	Heptacos-1-ene	378	C27H54	015306-27-1	87
4	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	86
5	Cyclodocosane, ethyl-	336	C24H48	1000151-22-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFW8

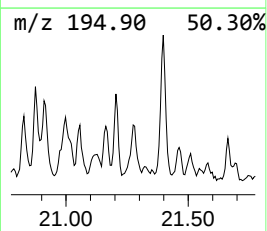
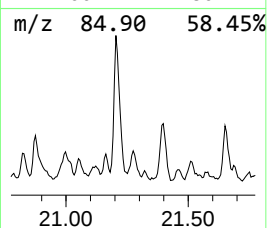
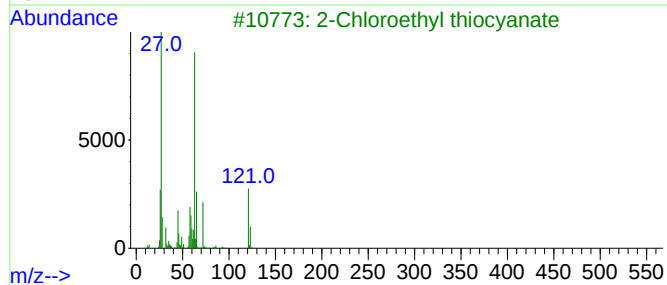
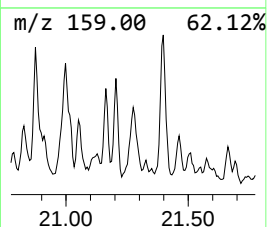
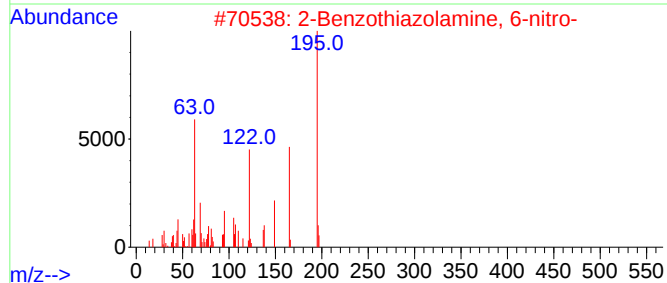
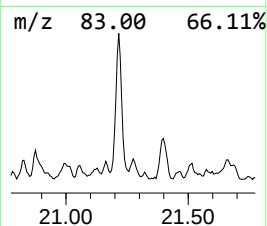
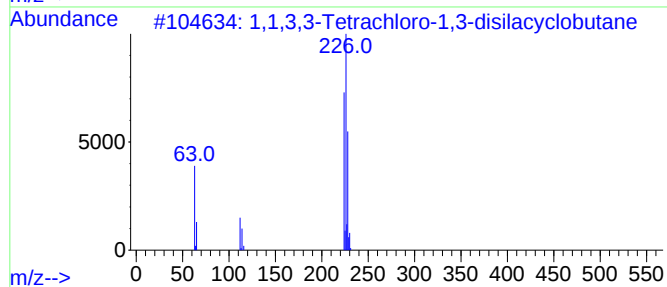
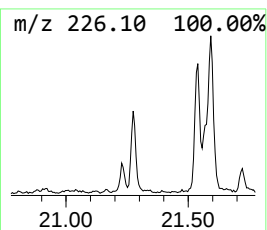
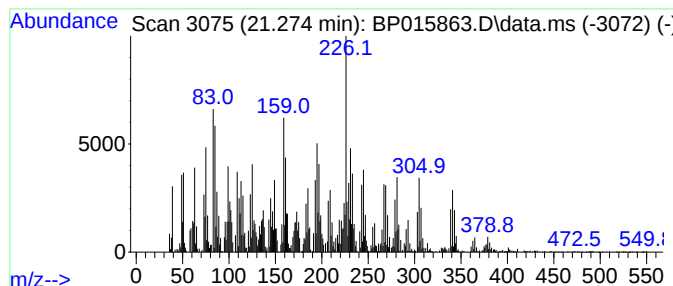
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Quant Title : SVOA CALIBRATION

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

Peak Number 28 unknown-13 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.275	8.20 ng/ul	923522	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,3,3-Tetrachloro-1,3-disilacy...	224	C2H4Cl4Si2	002146-97-6	38		
2	2-Benzothiazolamine, 6-nitro-	195	C7H5N3O2S	006285-57-0	14		
3	2-Chloroethyl thiocyanate	121	C3H4ClNS	1000342-25-2	14		
4	Benzene, 1,2,3-trichloro-4-methyl-	194	C7H5Cl3	007359-72-0	12		
5	2,6-Dichlorobenzylsulfonyl chloride	258	C7H5Cl3O2S	085952-31-4	12		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

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ClientSampleId :
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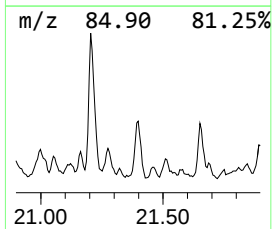
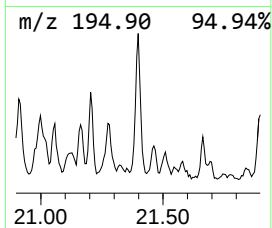
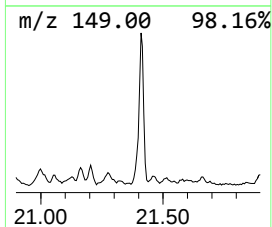
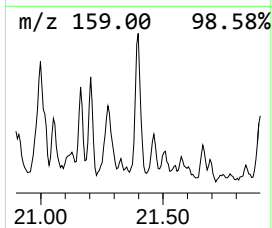
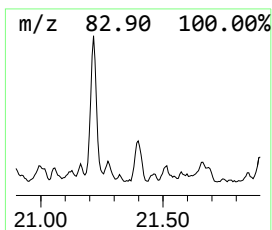
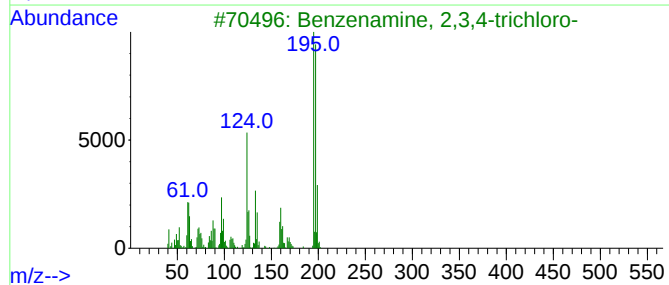
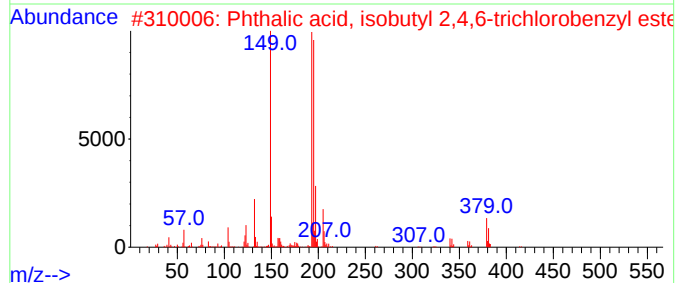
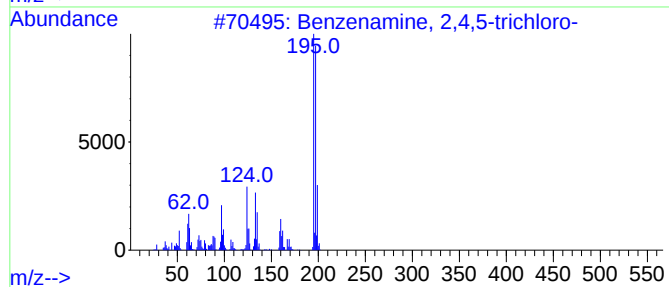
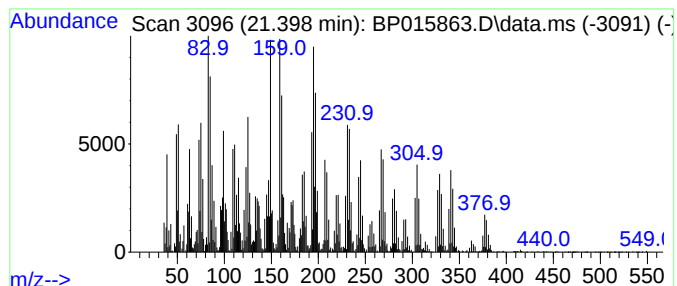
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Quant Title : SVOA CALIBRATION

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

Peak Number 29 unknown-14 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.398	21.76 ng/ul	2449030	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	44
2			Phthalic acid, isobutyl 2,4,6-tr...	414	C19H17Cl3O4	1000382-87-7	44
3			Benzenamine, 2,3,4-trichloro-	195	C6H4Cl3N	000634-67-3	44
4			Benzenamine, 2,4,5-trichloro-	195	C6H4Cl3N	000636-30-6	25
5			Selenolo[2,3-b]pyridine, 3-methyl-	197	C8H7NSe	055108-56-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
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Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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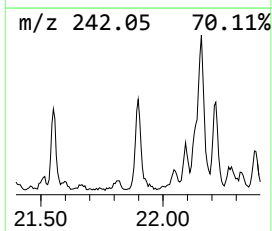
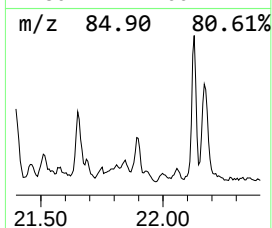
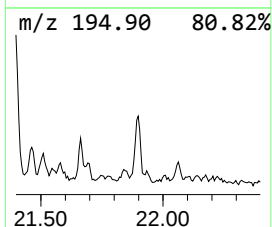
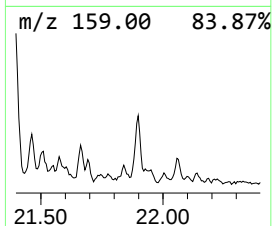
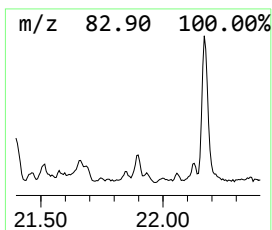
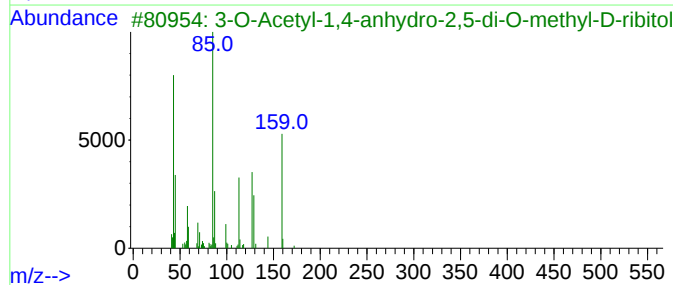
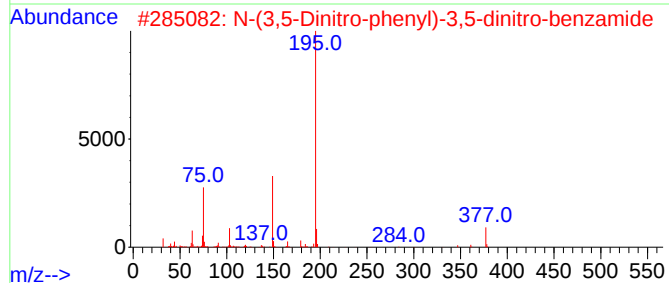
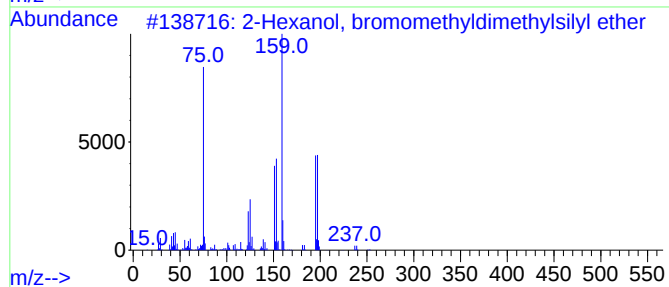
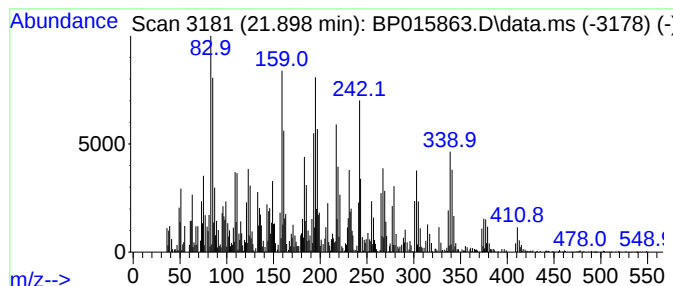
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 unknown-15 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.898	4.25 ng/ul	478360	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexanol, bromomethyldimethylsi...	252	C9H21BrOSi	1000375-64-8	30		
2	N-(3,5-Dinitro-phenyl)-3,5-dinit...	377	C13H7N5O9	058338-55-9	25		
3	3-O-Acetyl-1,4-anhydro-2,5-di-O-...	204	C9H16O5	1000357-22-8	12		
4	2,2',4',5'-Tetrachloroacetanilide	271	C8H5Cl4NO	023595-42-8	12		
5	3-Aminoxy-4-chlorobutyric acid,...	181	C6H12ClNO3	1000186-78-9	12		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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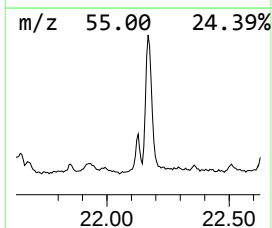
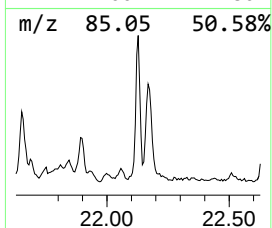
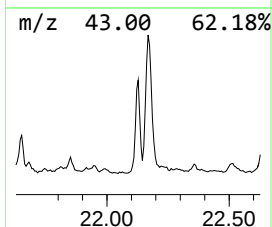
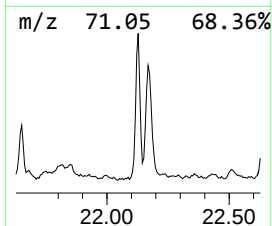
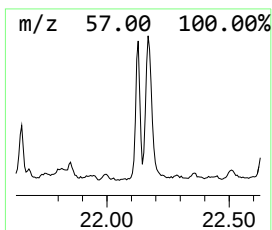
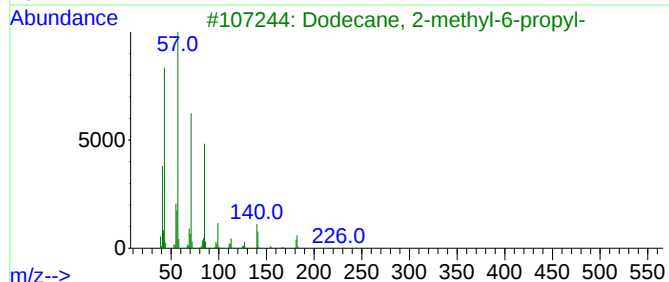
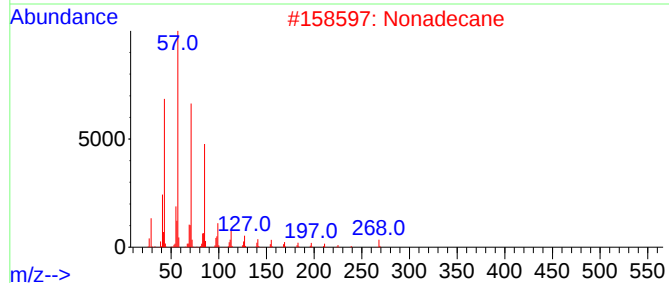
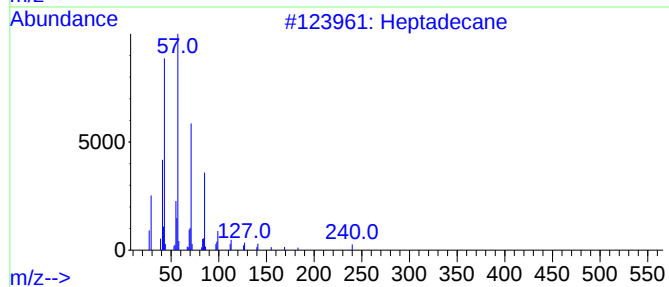
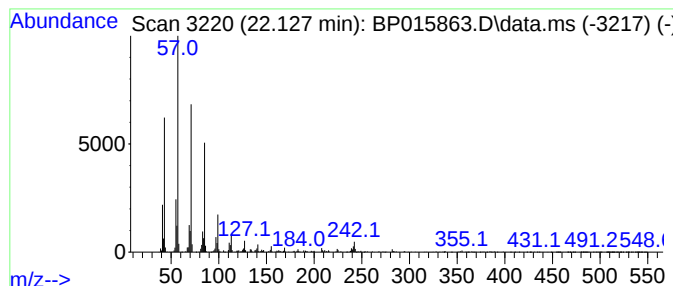
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.127	5.34 ng/ul	601140	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	94
2			Nonadecane	268	C19H40	000629-92-5	93
3			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	87
4			Heneicosane	296	C21H44	000629-94-7	87
5			Pentadecane	212	C15H32	000629-62-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
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Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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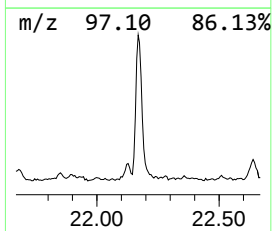
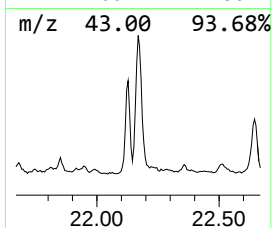
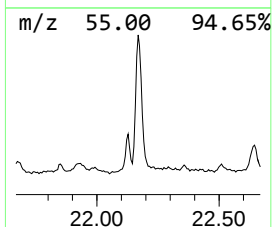
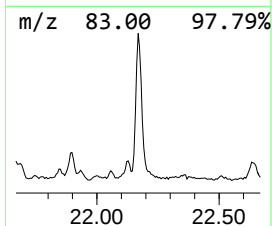
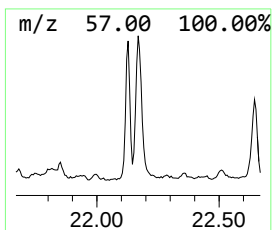
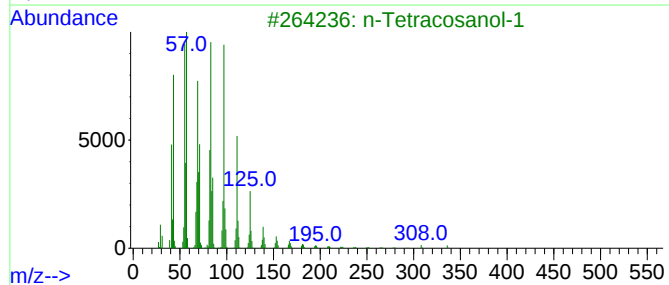
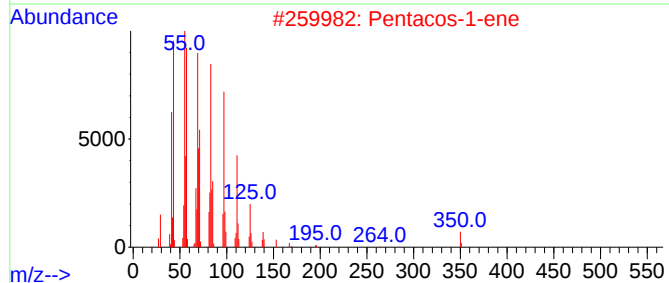
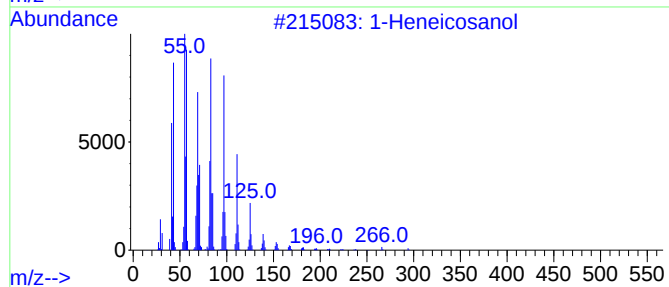
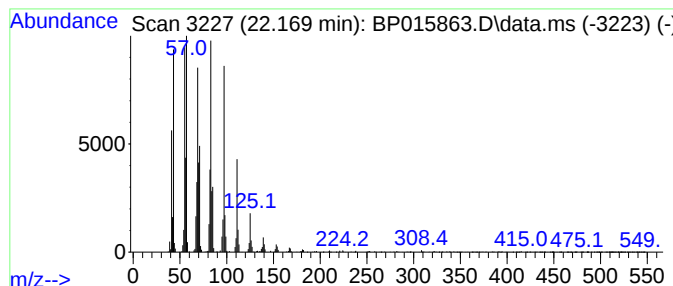
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 1-Heneicosanol Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.169	21.41 ng/ul	2410550	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosanol	312	C21H44O	015594-90-8	95
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			Behenic alcohol	326	C22H46O	000661-19-8	94
5			1-Octadecanol	270	C18H38O	000112-92-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
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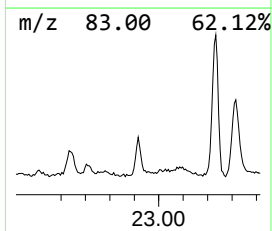
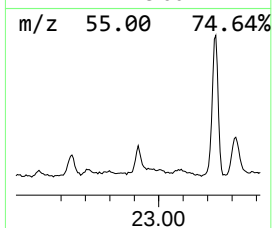
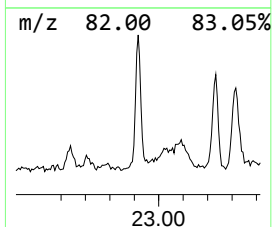
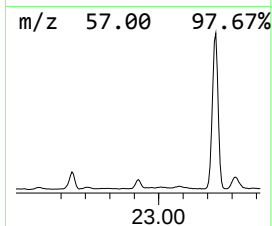
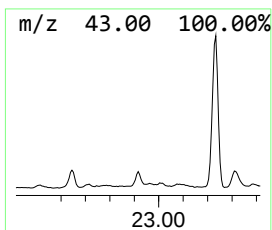
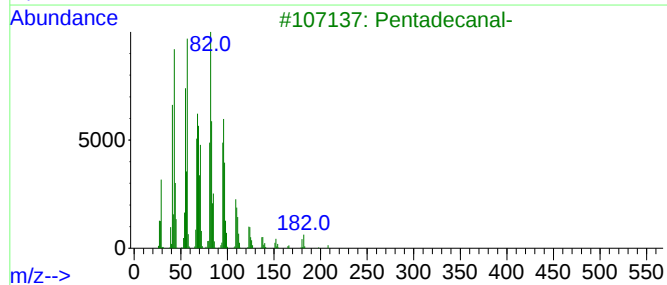
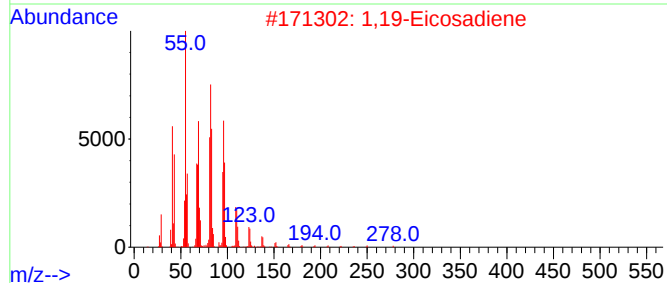
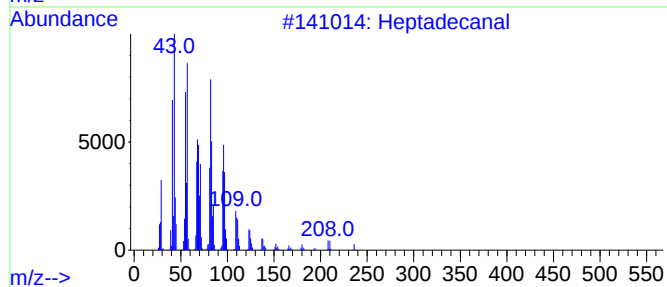
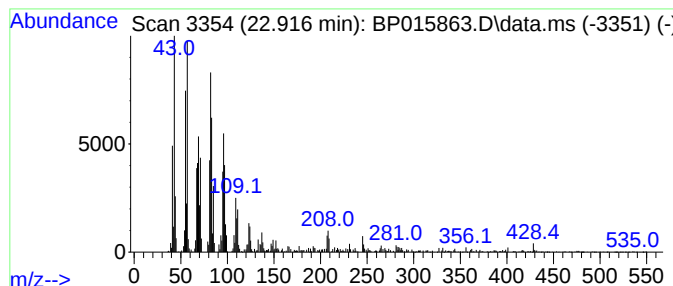
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Quant Title : SVOA CALIBRATION

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

Peak Number 33 Heptadecanal Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.916	3.17 ng/ul	508419	Perylene-d12	24.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecanal	254	C17H34O	1000376-70-0	93
2			1,19-Eicosadiene	278	C20H38	014811-95-1	93
3			Pentadecanal-	226	C15H30O	002765-11-9	92
4			Octadecanal	268	C18H36O	000638-66-4	91
5			Hexacosanal	380	C26H52O	026627-85-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
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Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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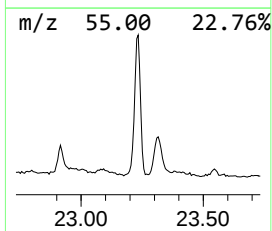
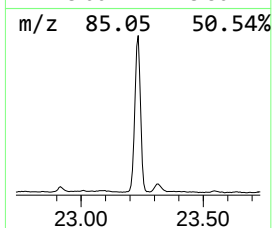
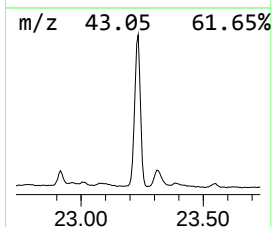
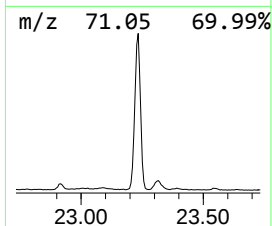
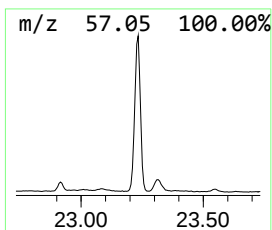
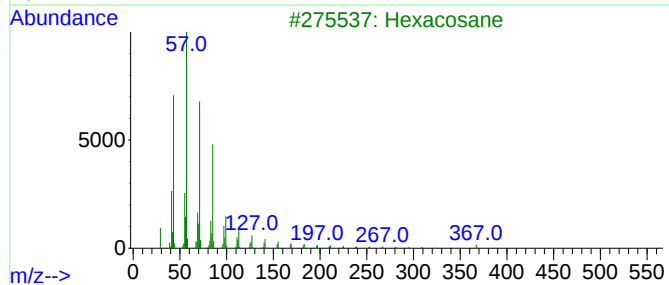
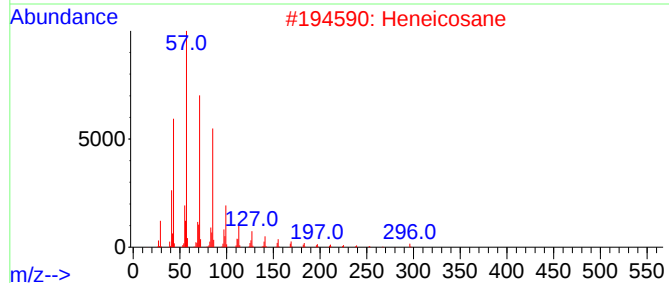
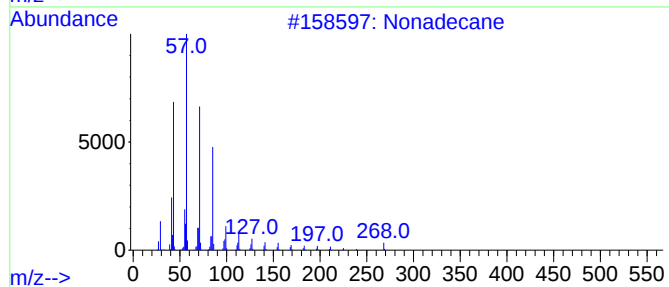
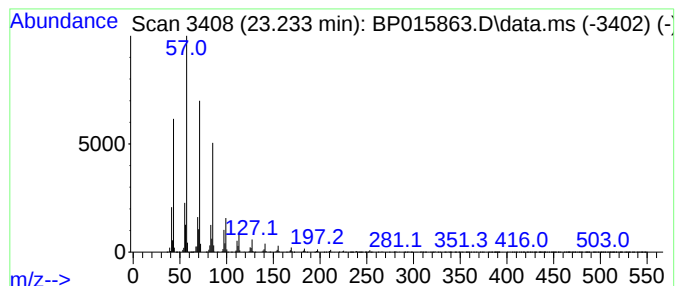
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.233	31.75 ng/ul	5098570	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonadecane	268	C19H40	000629-92-5	94
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
5		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFw8

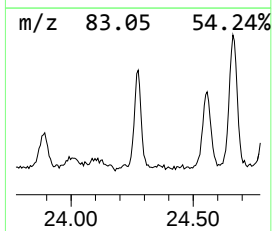
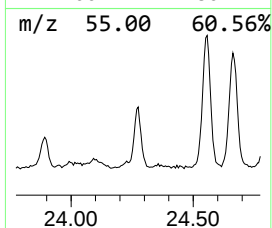
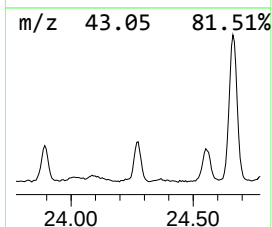
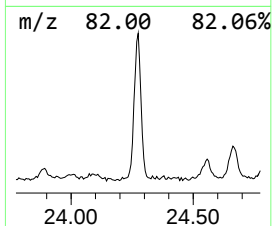
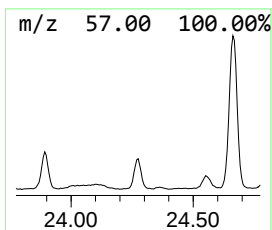
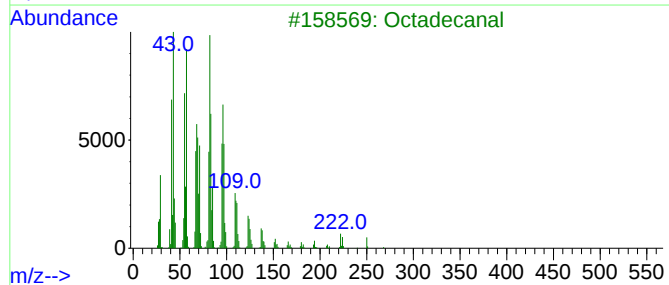
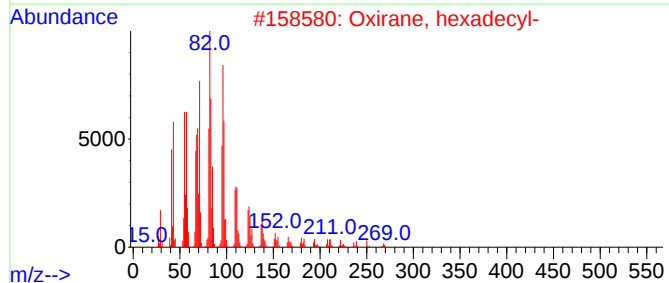
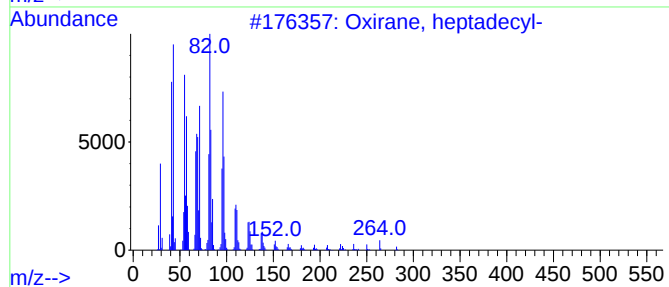
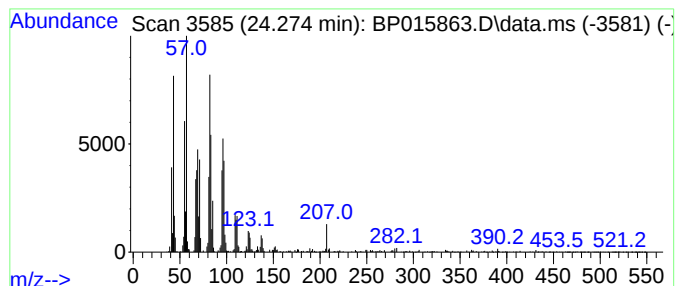
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Oxirane, heptadecyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.274	7.51 ng/ul	1205670	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Oxirane, heptadecyl-	282	C19H38O	067860-04-2	94
2		Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
3		Octadecanal	268	C18H36O	000638-66-4	91
4		Tetracosanal	352	C24H48O	057866-08-7	91
5		Octacosanal	408	C28H56O	022725-64-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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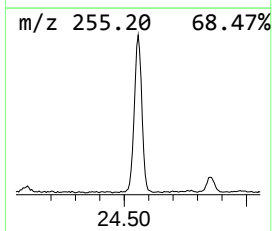
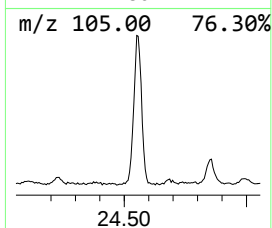
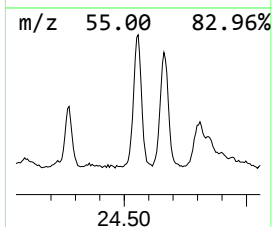
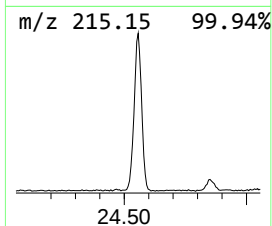
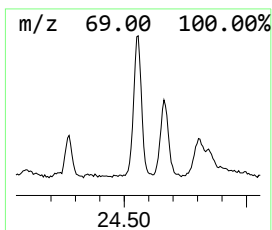
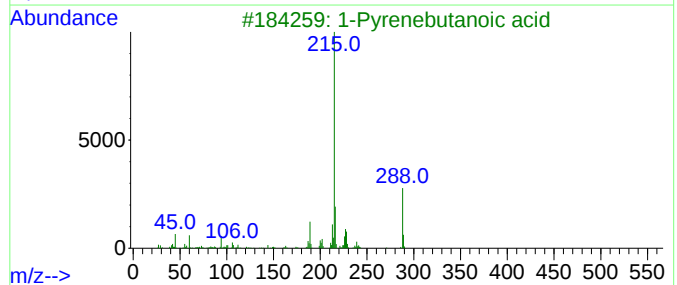
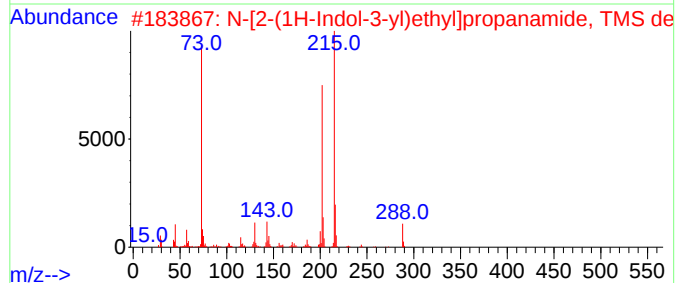
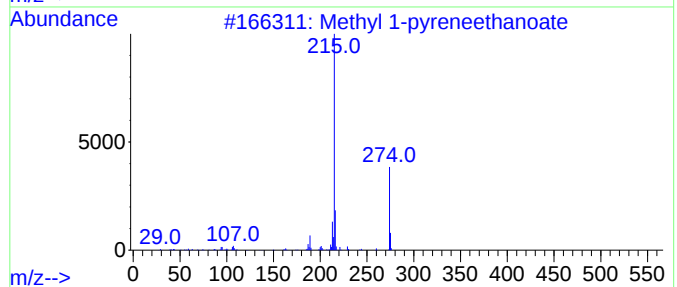
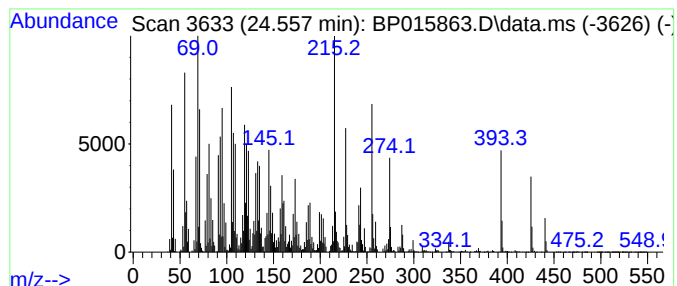
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 unknown-16 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.557	41.28 ng/ul	6627760	Perylene-d12	24.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	35
2			N-[2-(1H-Indol-3-yl)ethyl]propan...	288	C16H24N2OSi	1000493-19-0	25
3			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	15
4			Indanidine	215	C11H13N5	085392-79-6	15
5			10-Fluoro-2,4-dimethylpyrimido[1...	215	C12H10FN3	1000388-01-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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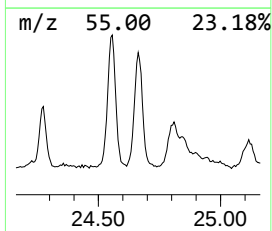
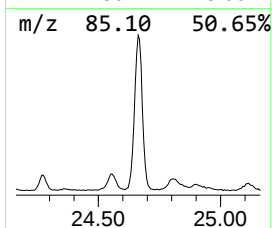
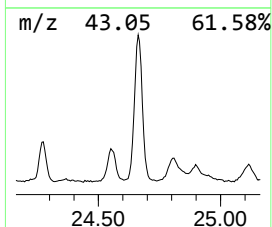
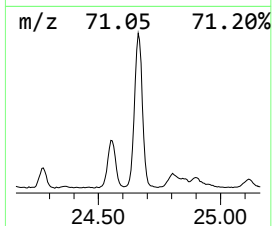
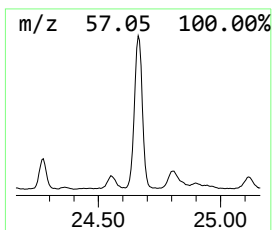
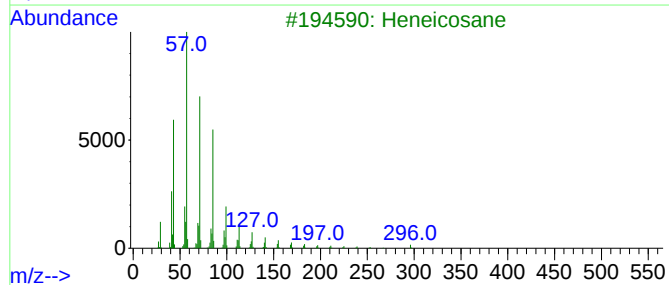
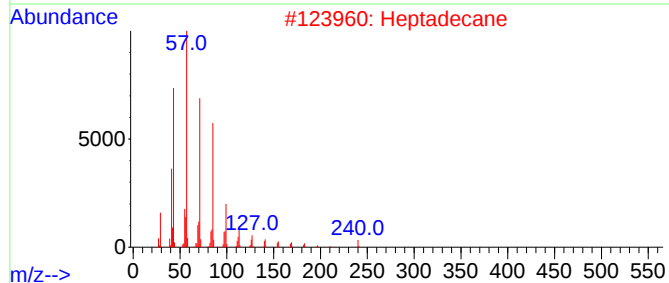
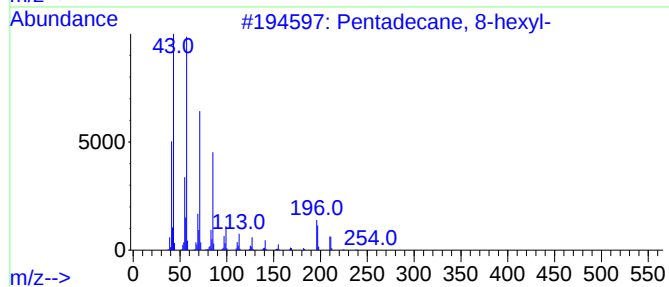
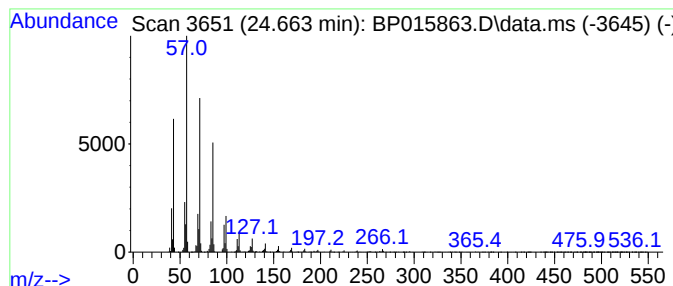
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.663	23.65 ng/ul	3798290	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Heptadecane	240	C17H36	000629-78-7	93
3		Heneicosane	296	C21H44	000629-94-7	91
4		9-Methylheneicosane	310	C22H46	070475-51-3	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFW8

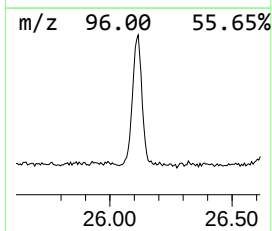
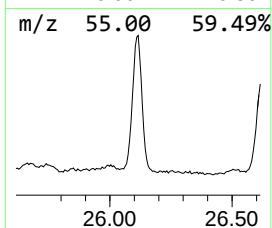
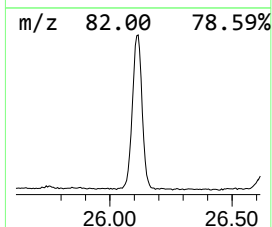
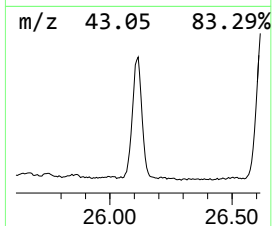
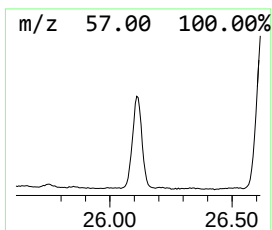
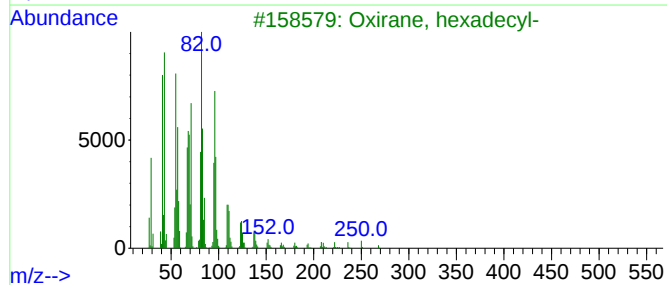
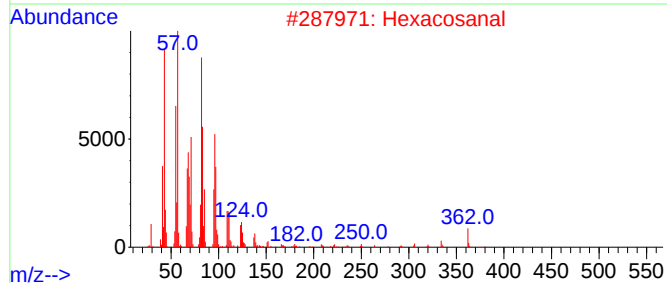
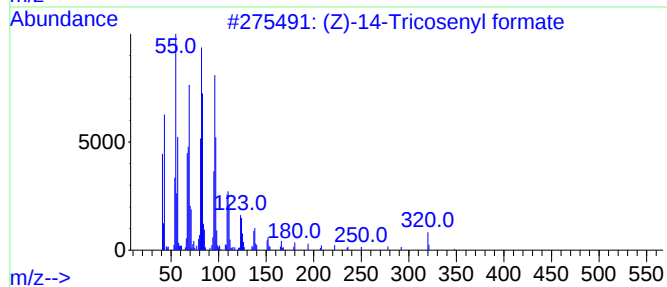
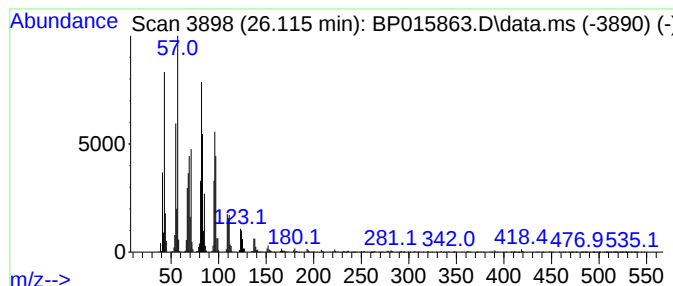
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (Z)-14-Tricosenyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.115	31.02 ng/ul	4981100	Perylene-d12	24.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(Z)-14-Tricosenyl formate			366	C24H46O2	077899-10-6	95
2	Hexacosanal			380	C26H52O	026627-85-0	91
3	Oxirane, hexadecyl-			268	C18H36O	007390-81-0	91
4	Octadecanal			268	C18H36O	000638-66-4	90
5	Octadecanal			268	C18H36O	000638-66-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015863.D
Acq On : 01 Jul 2023 01:54
Operator : MA/JU
Sample : 03161-12
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFw8

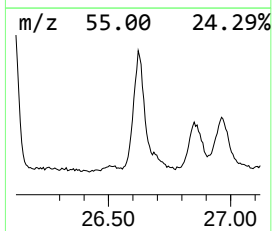
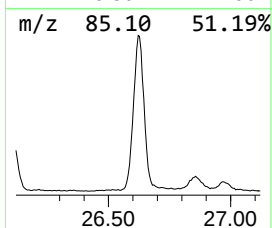
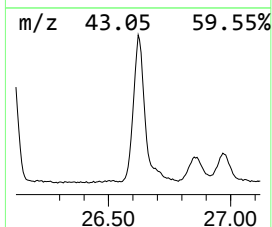
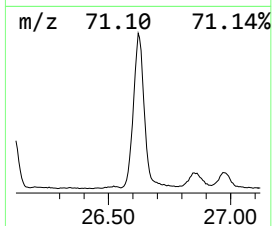
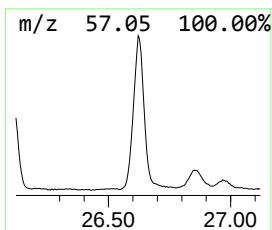
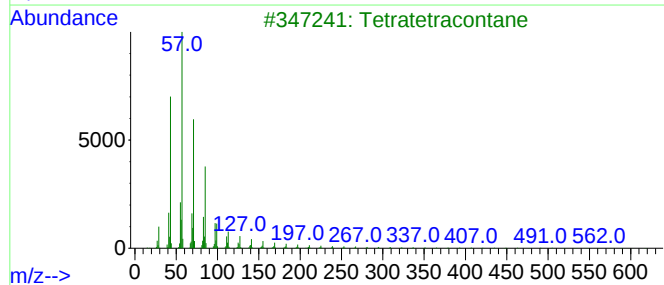
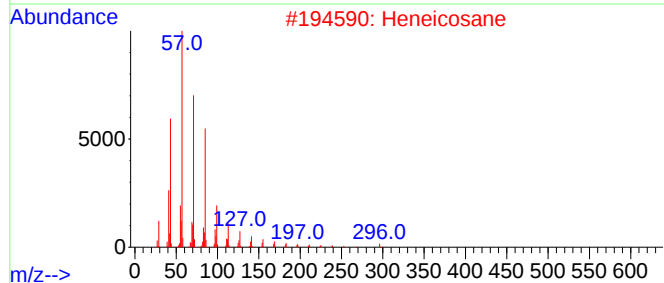
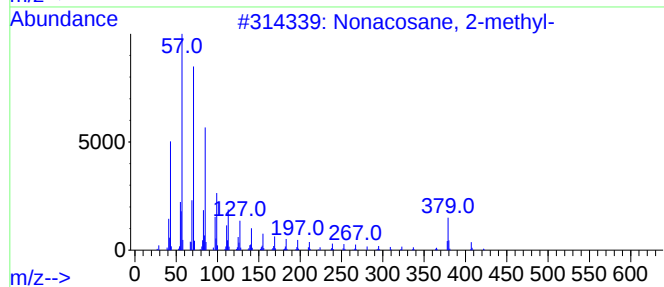
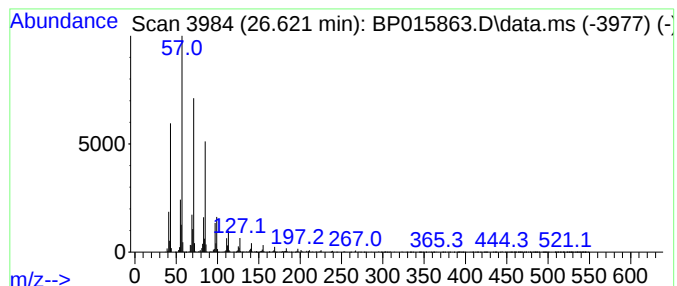
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.621	36.01 ng/ul	5783090	Perylene-d12	24.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosane, 2-methyl-	422	C30H62	001560-75-4	94
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetratetracontane	619	C44H90	007098-22-8	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Octacosane, 1-iodo-	520	C28H57I	1000406-32-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015863.D
 Acq On : 01 Jul 2023 01:54
 Operator : MA/JU
 Sample : 03161-12
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFW8

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
2H-Pyran, tetra...	3.922	5.0	ng/ul	400583	1	8.052	1585950	20.0
Benzophenone	16.069	6.6	ng/ul	993609	3	14.704	3005890	20.0
Pentadecanoic acid	17.334	2.5	ng/ul	359535	4	17.463	2930290	20.0
(E)-Hexadec-9-e...	18.263	9.0	ng/ul	1316120	4	17.463	2930290	20.0
n-Hexadecanoic ...	18.316	27.9	ng/ul	4093530	4	17.463	2930290	20.0
(DEL) Alkane: S...	19.192	2.9	ng/ul	419137	4	17.463	2930290	20.0
9,12-Octadecadi...	19.404	3.0	ng/ul	440229	4	17.463	2930290	20.0
Octadecanoic acid	19.539	10.9	ng/ul	1229030	5	21.551	2251450	20.0
unknown-01	19.933	5.5	ng/ul	618904	5	21.551	2251450	20.0
unknown-02	20.098	3.0	ng/ul	339501	5	21.551	2251450	20.0
unknown-03	20.133	5.6	ng/ul	625689	5	21.551	2251450	20.0
unknown-04	20.174	2.4	ng/ul	274259	5	21.551	2251450	20.0
unknown-05	20.227	4.8	ng/ul	535570	5	21.551	2251450	20.0
(DEL) Alkane: S...	20.269	8.8	ng/ul	996115	5	21.551	2251450	20.0
unknown-06	20.339	8.7	ng/ul	976217	5	21.551	2251450	20.0
unknown-07	20.410	5.4	ng/ul	609522	5	21.551	2251450	20.0
Benzenamine, 2,...	20.457	8.7	ng/ul	984361	5	21.551	2251450	20.0
Bornane, 2,2,5-...	20.575	15.0	ng/ul	1691400	5	21.551	2251450	20.0
unknown-08	20.633	5.0	ng/ul	566320	5	21.551	2251450	20.0
unknown-09	20.827	4.7	ng/ul	532995	5	21.551	2251450	20.0
unknown-10	20.875	18.4	ng/ul	2066670	5	21.551	2251450	20.0
2,2,5-Endo,6-ex...	20.998	13.7	ng/ul	1537420	5	21.551	2251450	20.0
unknown-11	21.051	4.5	ng/ul	503431	5	21.551	2251450	20.0
unknown-12	21.163	5.8	ng/ul	650145	5	21.551	2251450	20.0
(DEL) Alkane: S...	21.210	28.8	ng/ul	3246300	5	21.551	2251450	20.0
unknown-13	21.275	8.2	ng/ul	923522	5	21.551	2251450	20.0
unknown-14	21.398	21.8	ng/ul	2449030	5	21.551	2251450	20.0
unknown-15	21.898	4.3	ng/ul	478360	5	21.551	2251450	20.0
(DEL) Alkane: S...	22.127	5.3	ng/ul	601140	5	21.551	2251450	20.0
1-Heneicosanol	22.169	21.4	ng/ul	2410550	5	21.551	2251450	20.0
Heptadecanal	22.916	3.2	ng/ul	508419	6	24.092	3211500	20.0
(DEL) Alkane: S...	23.233	31.8	ng/ul	5098570	6	24.092	3211500	20.0
Oxirane, heptad...	24.274	7.5	ng/ul	1205670	6	24.092	3211500	20.0
unknown-16	24.557	41.3	ng/ul	6627760	6	24.092	3211500	20.0
(DEL) Alkane: S...	24.663	23.6	ng/ul	3798290	6	24.092	3211500	20.0
(Z)-14-Tricosen...	26.115	31.0	ng/ul	4981100	6	24.092	3211500	20.0
(DEL) Alkane: S...	26.621	36.0	ng/ul	5783090	6	24.092	3211500	20.0