

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
 Data File : BP020858.D
 Acq On : 09 Jul 2024 10:25
 Operator : MA/JU
 Sample : P3035-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CW2

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-BP062524.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020858.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	42	47	55	rVB	51359	77583	1.09%	0.062%
2	3.540	90	94	107	rVB	46144	89175	1.26%	0.071%
3	3.775	129	134	141	rVV	100548	132414	1.87%	0.105%
4	4.528	258	262	269	rVB2	37945	56355	0.79%	0.045%
5	4.881	317	322	332	rVB	82638	143504	2.02%	0.114%
6	5.834	478	484	498	rVB4	32451	73403	1.03%	0.058%
7	6.916	658	668	684	rVV	1149278	1882853	26.52%	1.493%
8	7.069	687	694	710	rVB	866368	1402199	19.75%	1.112%
9	7.269	720	728	736	rBV	1253047	1885281	26.55%	1.495%
10	7.340	736	740	750	rVB4	17029	50518	0.71%	0.040%
11	7.722	792	805	813	rBV	747878	1204716	16.97%	0.955%
12	8.216	883	889	894	rBV	28805	52103	0.73%	0.041%
13	8.434	917	926	940	rBV	1199946	2147772	30.25%	1.703%
14	8.546	940	945	953	rVB2	46905	84053	1.18%	0.067%
15	8.881	994	1002	1012	rBV	1054696	1830166	25.78%	1.451%
16	9.428	1089	1095	1102	rBV	67806	102999	1.45%	0.082%
17	9.587	1113	1122	1136	rBV	879972	1628265	22.93%	1.291%
18	10.128	1206	1214	1233	rBV	1388301	2460311	34.65%	1.951%
19	10.493	1266	1276	1282	rBV	1051498	1803607	25.40%	1.430%
20	10.645	1294	1302	1319	rBV	255553	666949	9.39%	0.529%
21	11.040	1361	1369	1377	rBV6	24167	57000	0.80%	0.045%
22	11.240	1396	1403	1410	rBV	80271	173355	2.44%	0.137%
23	13.287	1745	1751	1758	rBV3	18998	47895	0.67%	0.038%
24	13.351	1758	1762	1775	rVB2	40628	83165	1.17%	0.066%
25	13.528	1785	1792	1802	rBV3	31468	92106	1.30%	0.073%
26	13.663	1810	1815	1821	rBV4	28629	49025	0.69%	0.039%
27	13.757	1825	1831	1839	rVV	2555791	3414209	48.09%	2.707%
28	14.045	1869	1880	1897	rBV	2980810	4430675	62.41%	3.513%
29	14.181	1897	1903	1908	rVV	71622	112171	1.58%	0.089%
30	14.357	1922	1933	1939	rBV	1702821	2462356	34.68%	1.953%
31	14.416	1939	1943	1949	rVV	160133	238930	3.37%	0.189%
32	14.469	1949	1952	1963	rVB5	64151	102615	1.45%	0.081%
33	14.598	1963	1974	1994	rBV	1040053	2250315	31.70%	1.784%
34	14.751	1994	2000	2004	rVV2	96920	205688	2.90%	0.163%
35	14.816	2008	2011	2020	rVB2	116529	163855	2.31%	0.130%
36	15.163	2063	2070	2074	rBV5	19780	46461	0.65%	0.037%

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Integration Parameters: LSCINT.P

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	15.375	2098	2106	2112	rVV	3245402	5291713	74.53%	4.196%
38	15.428	2112	2115	2123	rVB	59280	88121	1.24%	0.070%
39	15.510	2123	2129	2139	rBV	1000213	1589636	22.39%	1.261%
40	15.633	2146	2150	2157	rVB4	47336	83137	1.17%	0.066%
41	15.957	2199	2205	2211	rBV6	63141	99987	1.41%	0.079%
42	16.110	2226	2231	2232	rBV4	38856	49913	0.70%	0.040%
43	16.304	2260	2264	2270	rVV7	23142	47612	0.67%	0.038%
44	16.569	2303	2309	2314	rBV8	49270	108484	1.53%	0.086%
45	16.686	2322	2329	2333	rBV6	40668	87856	1.24%	0.070%
46	16.822	2348	2352	2359	rVB	52763	83028	1.17%	0.066%
47	16.928	2363	2370	2376	rVV7	58702	124156	1.75%	0.098%
48	16.986	2376	2380	2385	rVB	57828	84665	1.19%	0.067%
49	17.039	2385	2389	2392	rBV3	55677	87815	1.24%	0.070%
50	17.075	2392	2395	2403	rVB3	92826	164288	2.31%	0.130%
51	17.175	2403	2412	2416	rBV	2099010	3152429	44.40%	2.500%
52	17.216	2416	2419	2424	rVV	953903	1389975	19.58%	1.102%
53	17.280	2424	2430	2439	rVV2	3458011	5910772	83.25%	4.687%
54	17.351	2439	2442	2450	rVB5	117681	185094	2.61%	0.147%
55	17.604	2477	2485	2488	rBV2	123545	241027	3.39%	0.191%
56	17.639	2488	2491	2494	rVB3	42954	52316	0.74%	0.041%
57	17.804	2510	2519	2527	rBV	669330	1122014	15.80%	0.890%
58	17.916	2533	2538	2545	rVB5	114403	185364	2.61%	0.147%
59	17.992	2545	2551	2556	rBV6	79221	159208	2.24%	0.126%
60	18.057	2557	2562	2563	rBV2	180938	247699	3.49%	0.196%
61	18.075	2563	2565	2568	rVV	146194	183132	2.58%	0.145%
62	18.116	2568	2572	2580	rVB2	1528572	2080947	29.31%	1.650%
63	18.286	2595	2601	2606	rBV2	595571	1019290	14.36%	0.808%
64	18.345	2606	2611	2615	rVB2	322282	516864	7.28%	0.410%
65	18.392	2615	2619	2626	rBV2	338336	485238	6.83%	0.385%
66	18.574	2645	2650	2653	rBV	253971	393510	5.54%	0.312%
67	18.651	2660	2663	2667	rVB	130034	165537	2.33%	0.131%
68	18.727	2673	2676	2678	rBV2	66168	73667	1.04%	0.058%
69	18.763	2678	2682	2687	rVB	205296	272714	3.84%	0.216%
70	18.863	2696	2699	2704	rVB3	88701	120672	1.70%	0.096%
71	19.033	2722	2728	2731	rBV3	109234	225962	3.18%	0.179%
72	19.080	2732	2736	2741	rVV3	324471	542271	7.64%	0.430%
73	19.139	2741	2746	2749	rVV	402185	675077	9.51%	0.535%
74	19.180	2749	2753	2761	rVB4	482867	892017	12.56%	0.707%
75	19.310	2768	2775	2783	rBV	2569178	4070841	57.34%	3.228%
76	19.451	2795	2799	2806	rVB3	708181	1043609	14.70%	0.828%

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Stop Thrs : 0

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77	19.527	2809	2812	2816	rVB3	135148	164142	2.31%	0.130%
78	19.657	2828	2834	2838	rBV2	4240142	7099776	100.00%	5.630%
79	19.686	2838	2839	2849	rVB	1826905	2034391	28.65%	1.613%
80	19.927	2876	2880	2886	rBV2	240809	380521	5.36%	0.302%
81	20.210	2924	2928	2931	rBV3	296190	398272	5.61%	0.316%
82	20.310	2942	2945	2949	rBV	191984	282408	3.98%	0.224%
83	20.757	3018	3021	3024	rVB2	221197	258862	3.65%	0.205%
84	20.998	3059	3062	3069	rVB	217387	360737	5.08%	0.286%
85	21.280	3105	3110	3117	rVB3	1737932	2841107	40.02%	2.253%
86	21.527	3149	3152	3158	rVB	526822	733218	10.33%	0.581%
87	21.615	3160	3167	3172	rVV2	2428482	5317334	74.89%	4.216%
88	21.663	3172	3175	3186	rVB	1307254	2252504	31.73%	1.786%
89	22.492	3311	3316	3319	rBV	1039264	1680683	23.67%	1.333%
90	22.533	3319	3323	3332	rVB3	1541122	2994653	42.18%	2.375%
91	23.598	3499	3504	3511	rVB	733994	1474379	20.77%	1.169%
92	23.909	3551	3557	3565	rBV	894415	2493015	35.11%	1.977%
93	24.068	3578	3584	3593	rVV	2598606	6448343	90.82%	5.113%
94	24.162	3593	3600	3612	rVB	1332946	3516736	49.53%	2.789%
95	24.739	3690	3698	3705	rBV2	1711912	4302609	60.60%	3.412%
96	24.968	3730	3737	3746	rVB2	1084422	2973765	41.89%	2.358%
97	25.627	3842	3849	3862	rVB2	1112976	3201878	45.10%	2.539%
98	26.286	3950	3961	3974	rVB	2363815	6509265	91.68%	5.162%
99	26.433	3979	3986	3987	rBV3	650960	1313827	18.51%	1.042%
100	31.150	4779	4788	4790	rBV3	733834	1974258	27.81%	1.566%

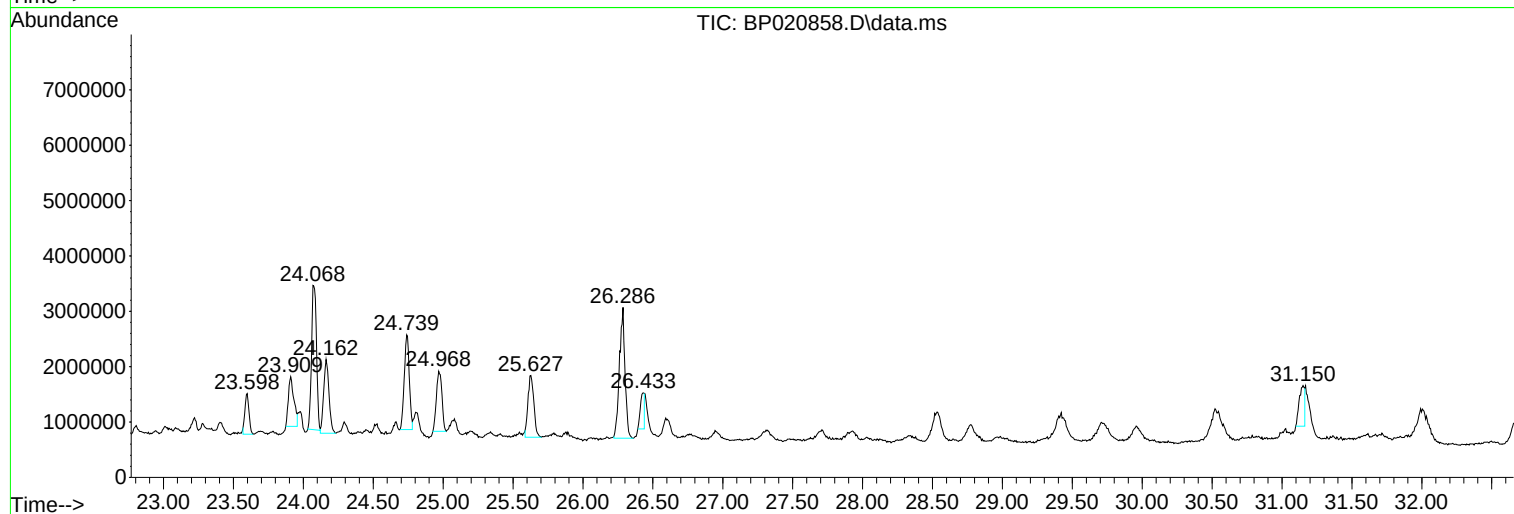
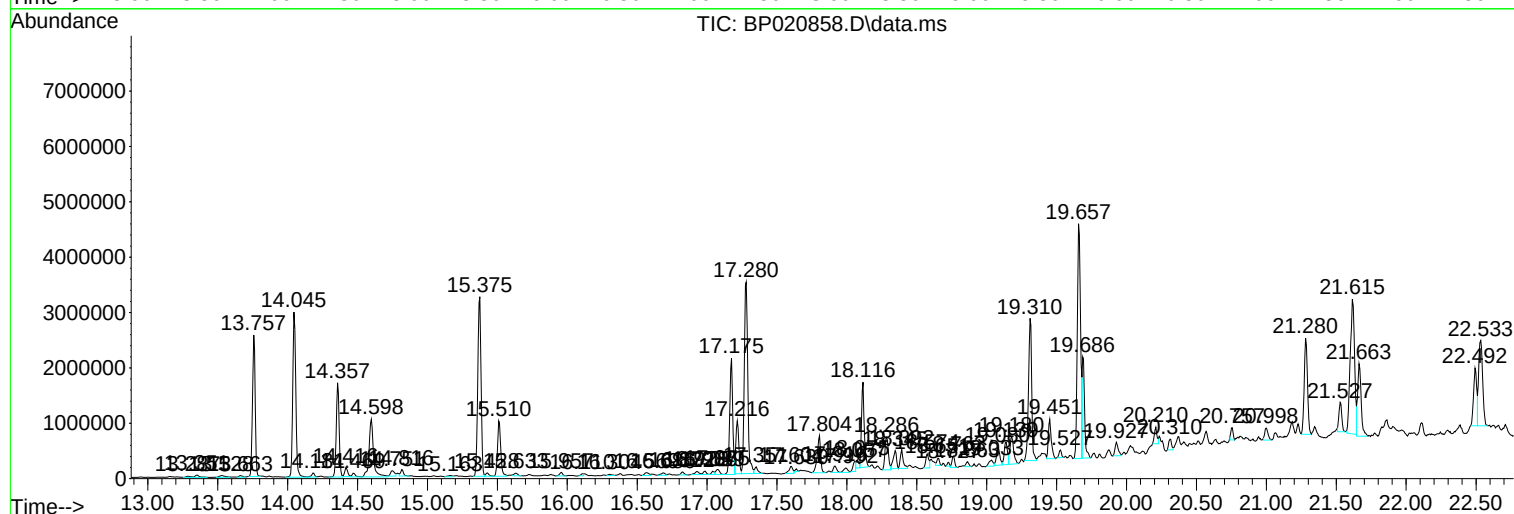
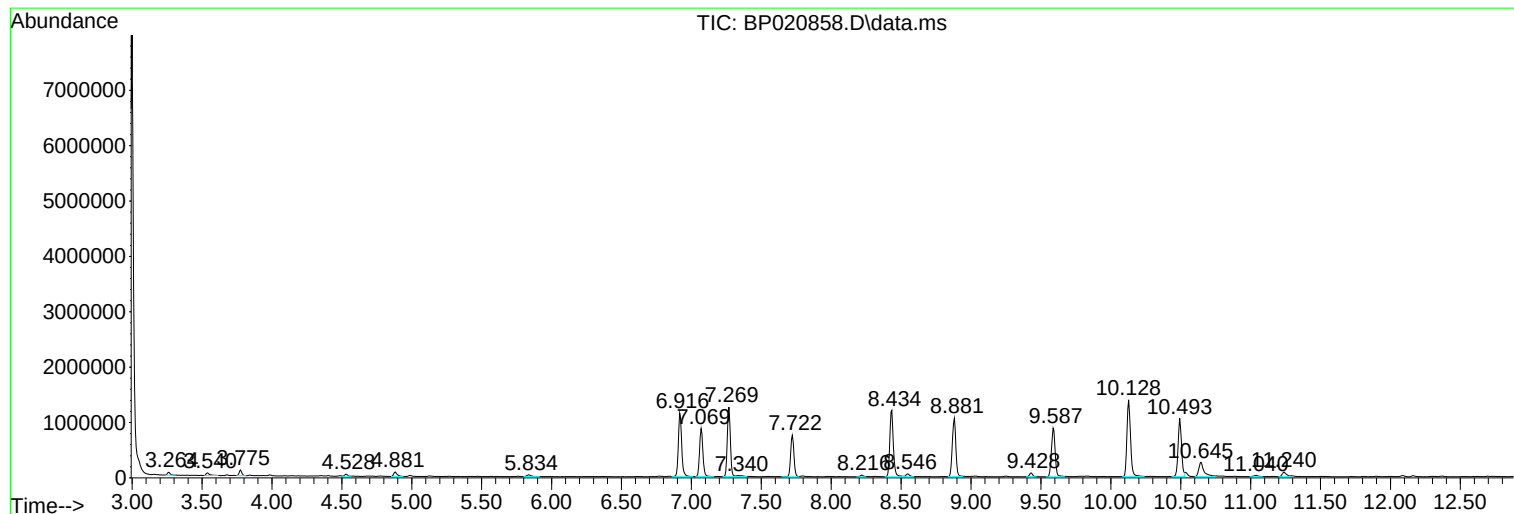
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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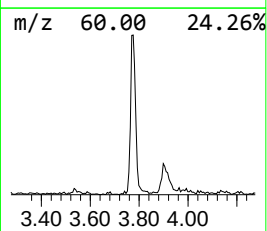
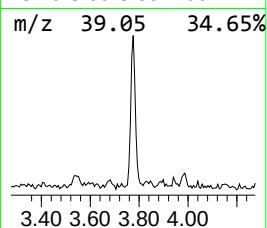
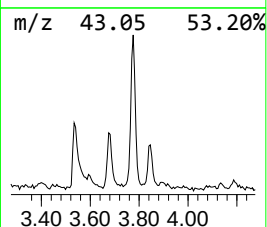
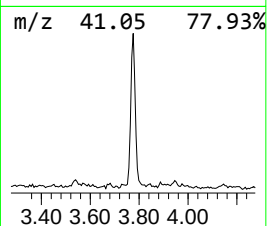
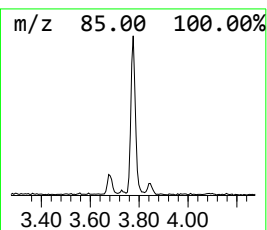
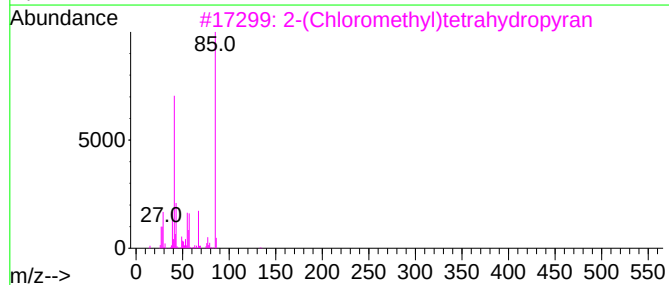
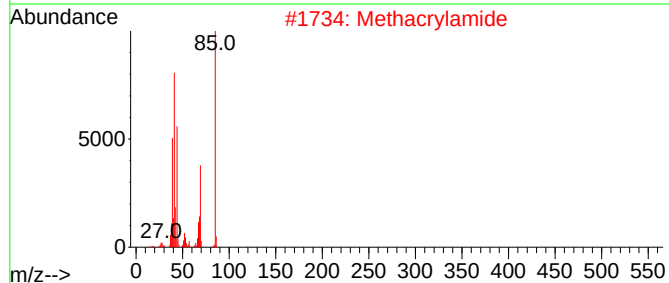
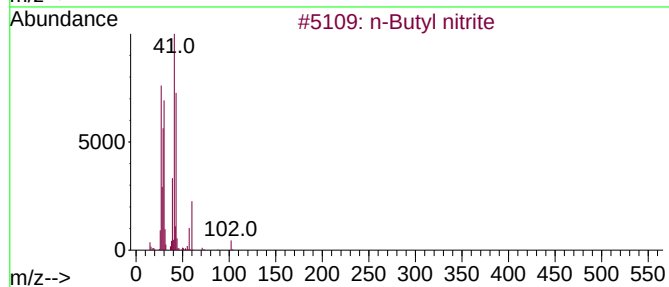
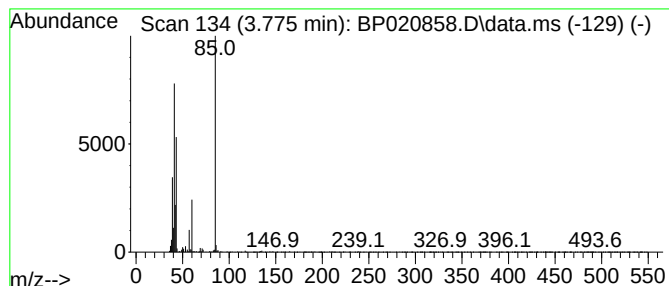
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.775	2.20 ng/ul	132414	1,4-Dichlorobenzene-d4	7.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Butyl nitrite	103	C4H9NO2	000544-16-1	14
2			Methacrylamide	85	C4H7NO	000079-39-0	12
3			2-(Chloromethyl)tetrahydropyran	134	C6H11ClO	018420-41-2	9
4			2,6-Dimethyl-8-(tetrahydropyran-...	254	C15H26O3	038290-53-8	9
5			Butanoic acid, 2-ethyl-2-methyl-...	144	C8H16O2	005296-70-8	9



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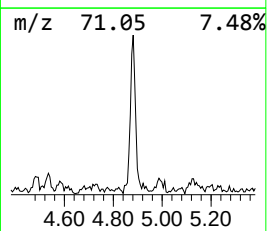
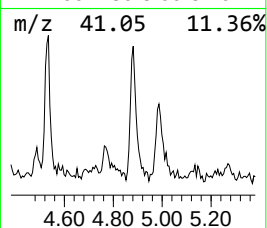
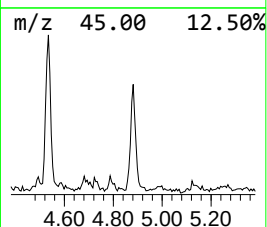
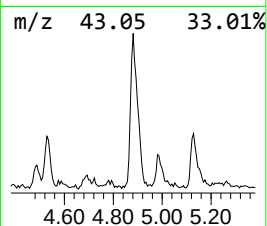
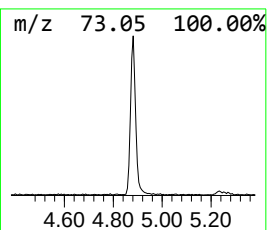
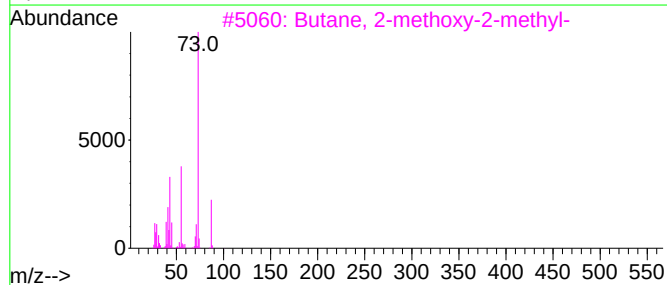
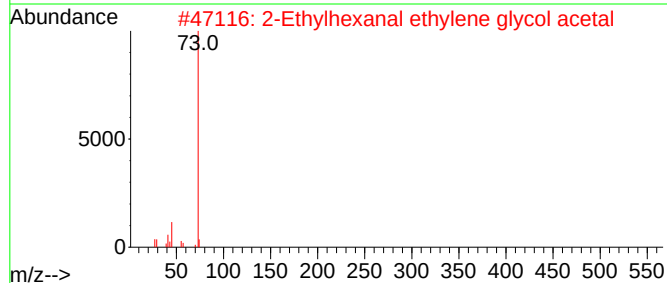
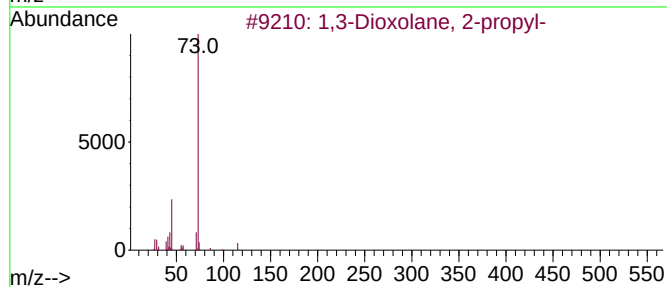
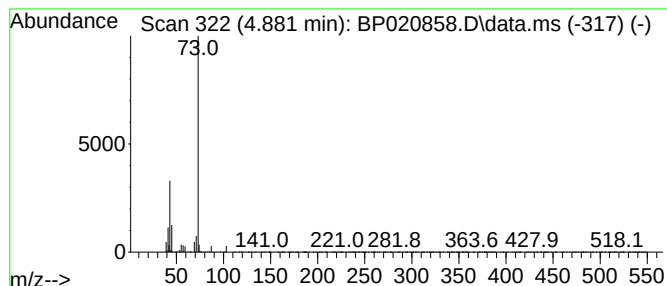
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-02 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.881	2.38 ng/ul	143504	1,4-Dichlorobenzene-d4	7.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	25
2			2-Ethylhexanal ethylene glycol a...	172	C10H20O2	1000431-01-2	9
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	9
4			1,3-Dioxolane, 2-heptyl-	172	C10H20O2	004359-57-3	9
5			2-(2-Bromoethyl)-1,3-dioxolane	180	C5H9BrO2	018742-02-4	9



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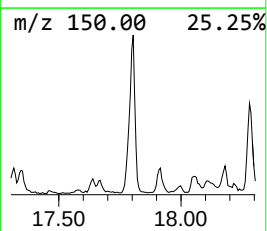
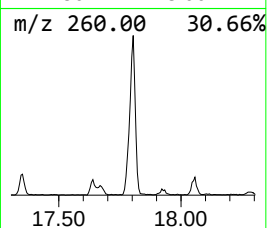
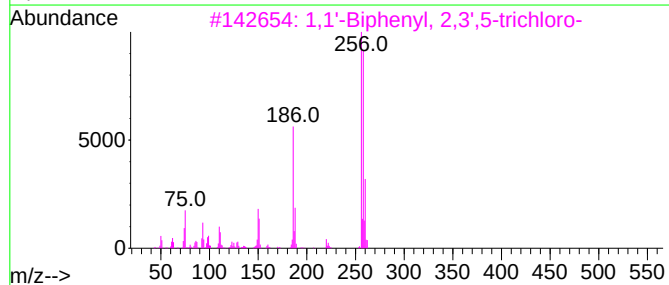
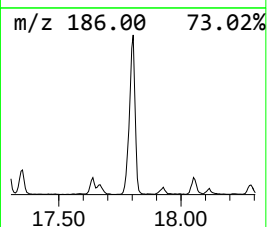
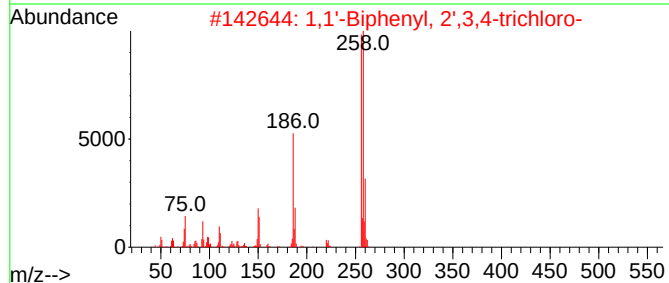
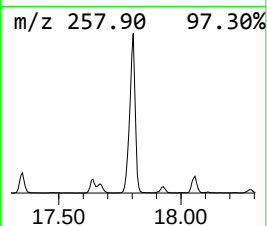
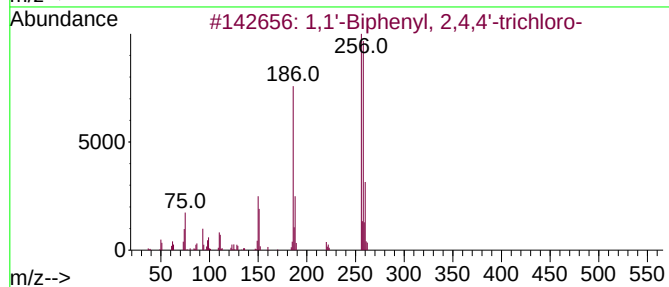
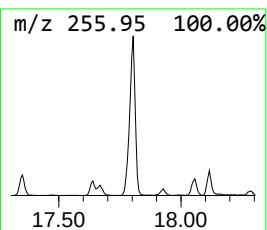
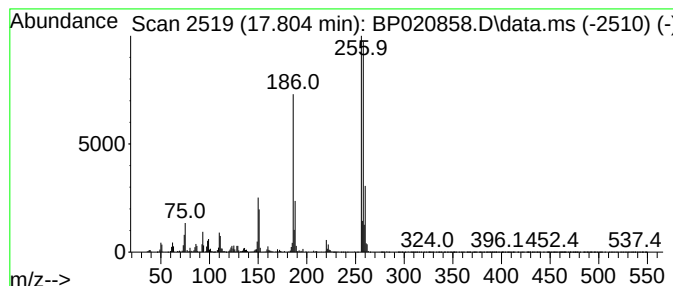
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1,1'-Biphenyl, 2,4,4'-trich... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.804	7.12 ng/ul	1122010	Phenanthrene-d10	17.174

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	99	
2	1,1'-Biphenyl, 2',3,4-trichloro-	256	C12H7Cl3	038444-86-9	99	
3	1,1'-Biphenyl, 2,3',5-trichloro-	256	C12H7Cl3	038444-81-4	99	
4	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99	
5	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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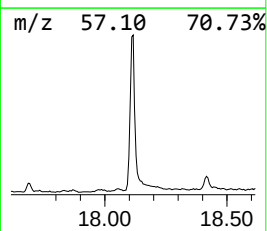
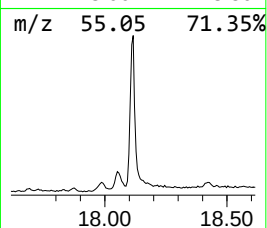
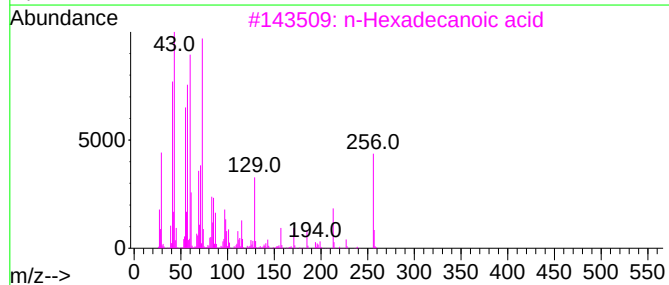
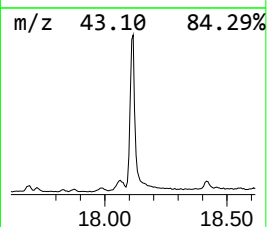
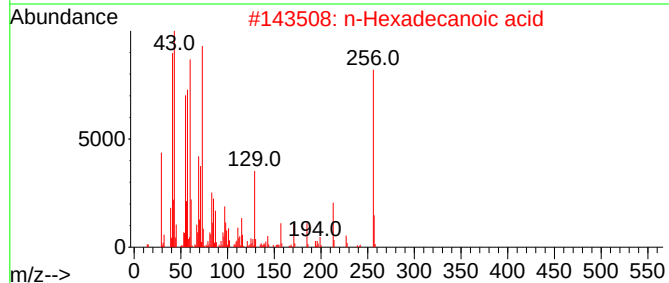
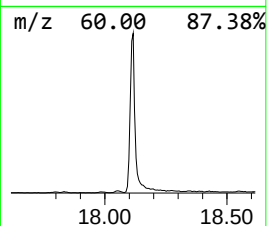
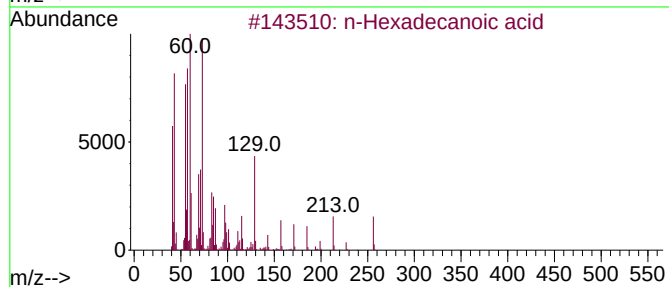
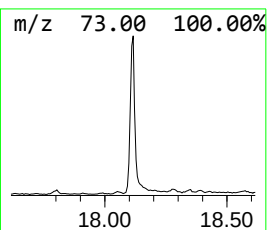
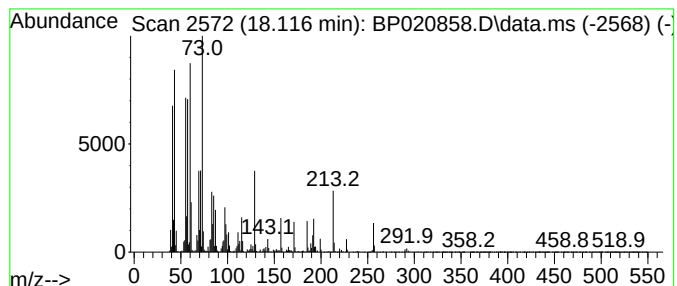
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	13.20 ng/ul	2080950	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
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Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

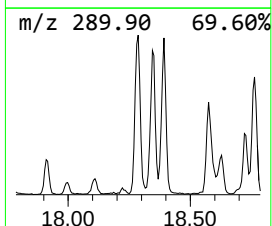
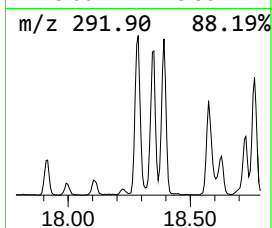
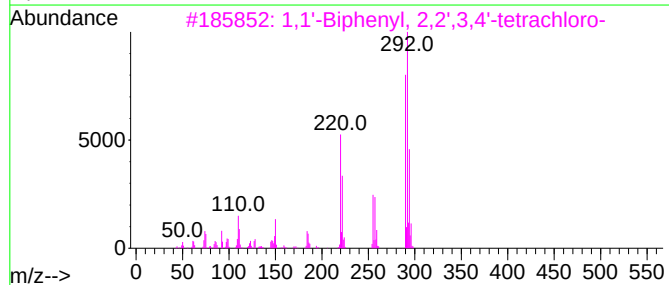
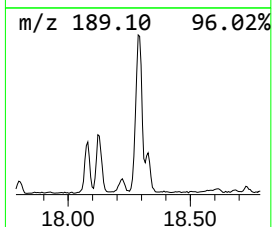
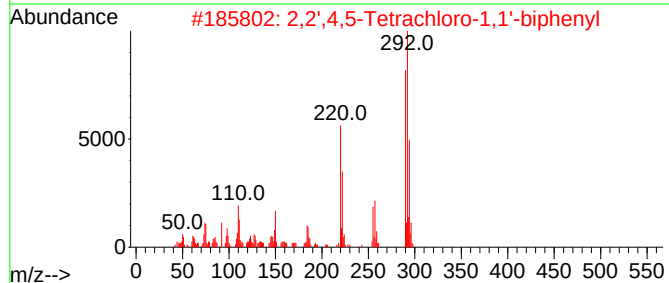
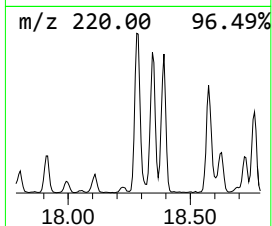
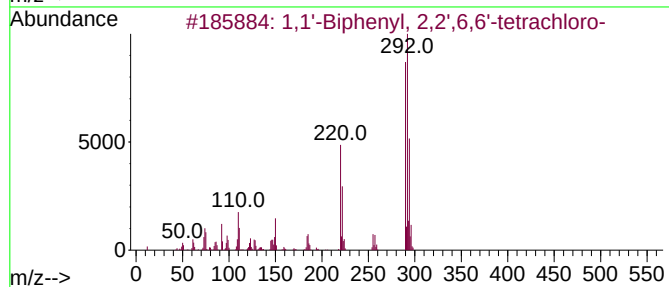
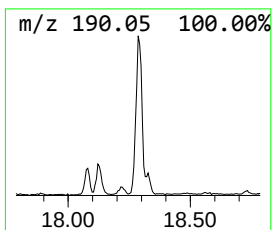
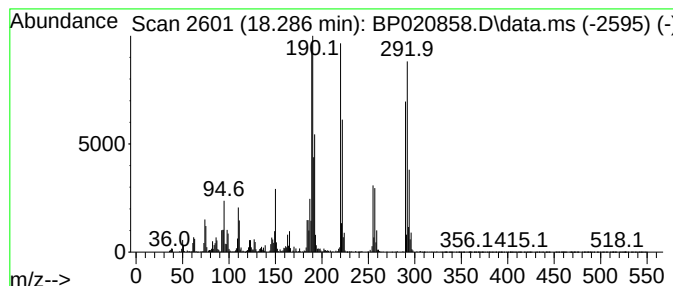
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2',6,6'-te... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.286	6.47 ng/ul	1019290	Phenanthrene-d10	17.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	97
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	97
3	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	97
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	97
5	2,3',4,5'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-52-7	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
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Operator : MA/JU
Sample : P3035-01
Misc :
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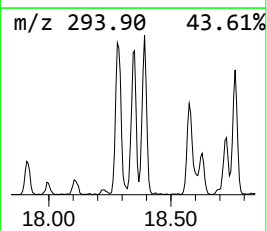
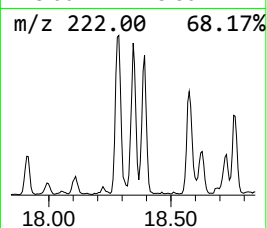
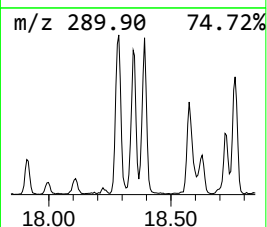
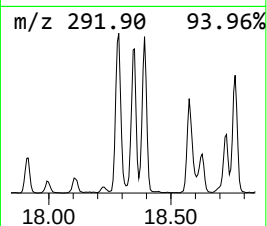
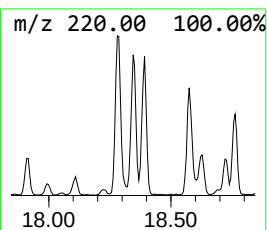
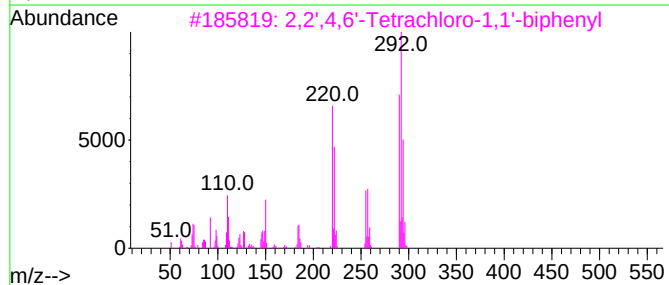
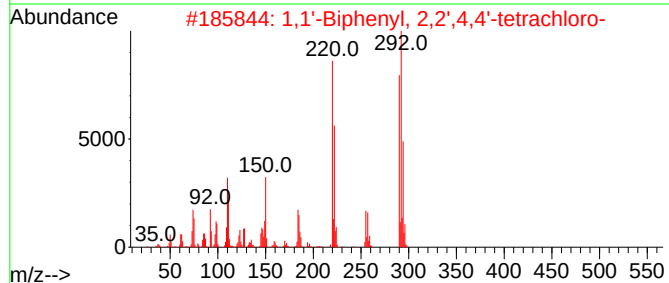
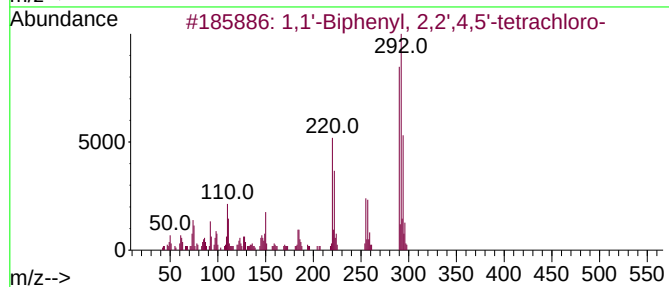
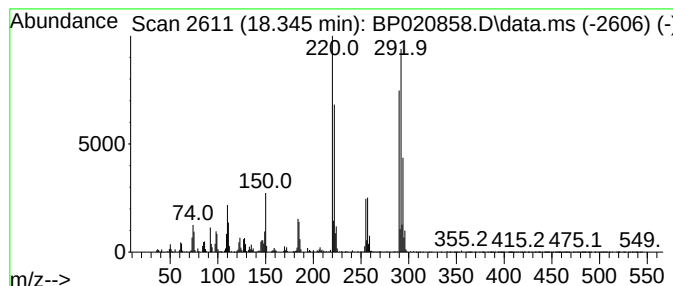
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 2,2',4,5'-te... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.345	3.28 ng/ul	516864	Phenanthrene-d10	17.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
3	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99
4	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
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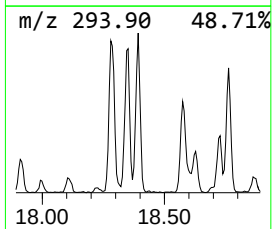
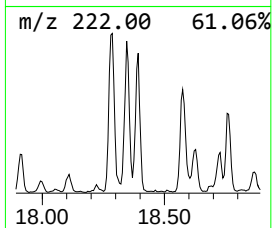
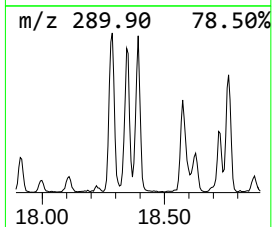
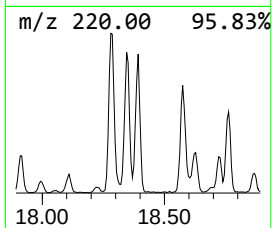
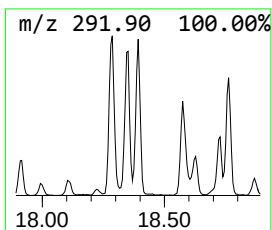
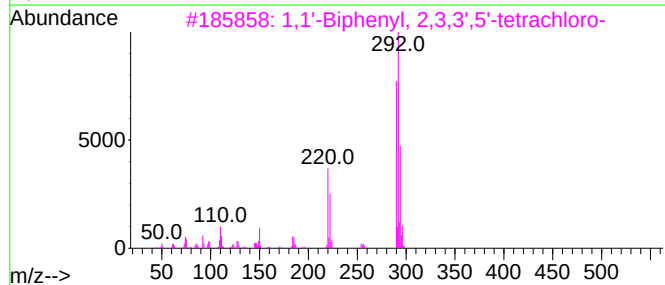
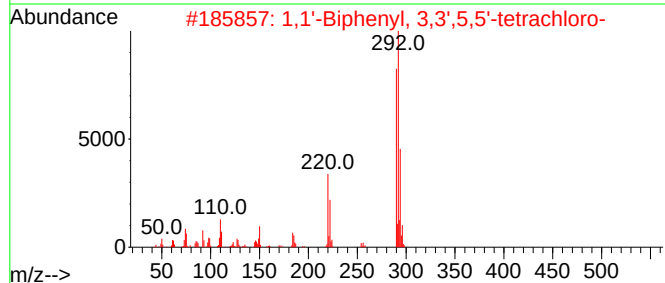
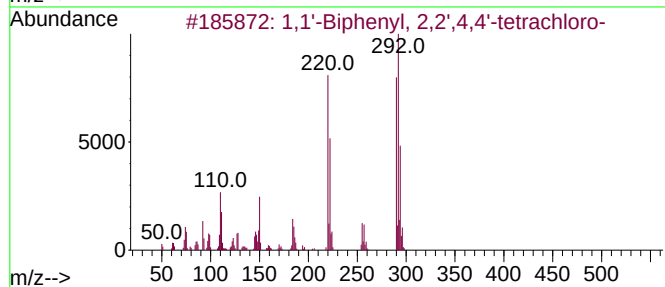
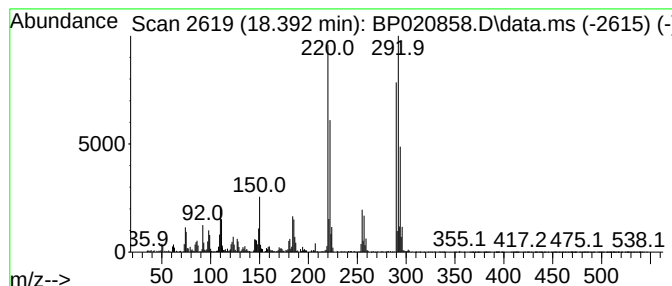
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.392	3.08 ng/ul	485238	Phenanthrene-d10	17.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99
2	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	99
3	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
4	2,3',4',5'-Tetrachloro-1,1'-biph...	290	C12H6Cl4	070362-48-0	99
5	1,1'-Biphenyl, 2,3',4',6'-tetrach...	290	C12H6Cl4	041464-46-4	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Misc :
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Instrument :
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ClientSampleId :
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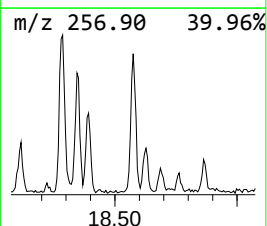
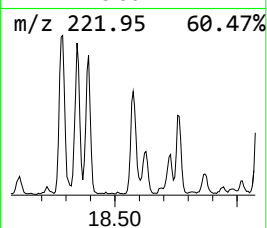
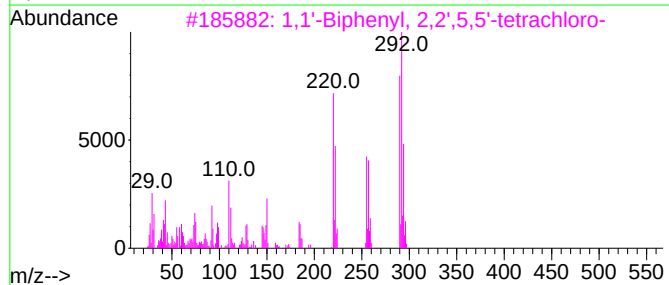
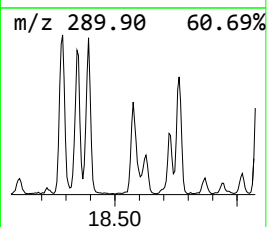
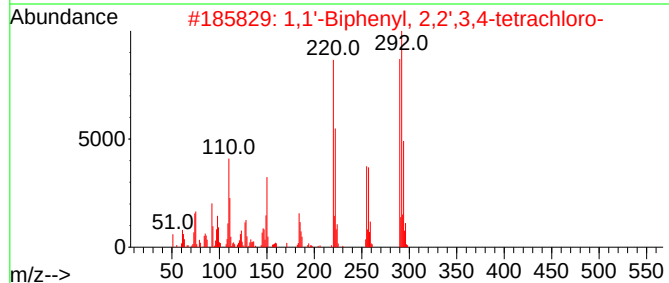
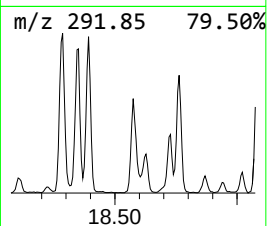
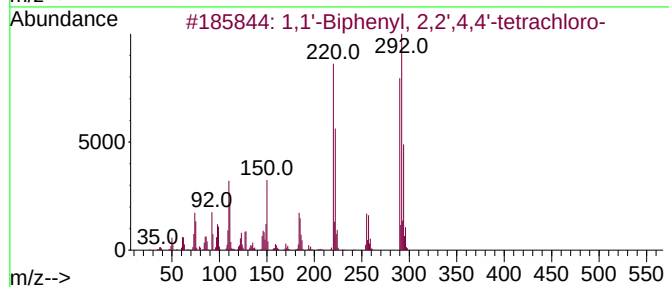
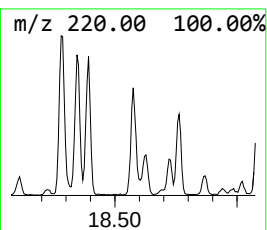
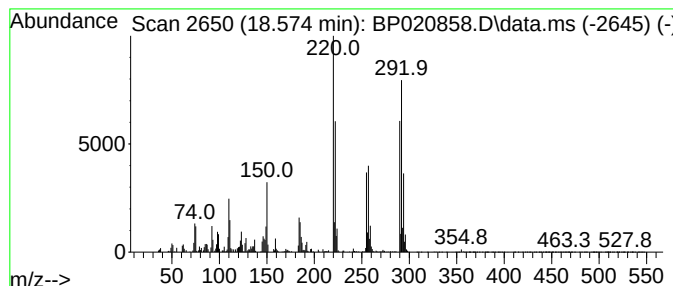
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,2',3,4-tet... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.575	2.50 ng/ul	393510	Phenanthrene-d10	17.174

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
2	1,1'-Biphenyl, 2,2',3,4-tetrachl...	290	C12H6Cl4	052663-59-9	99		
3	1,1'-Biphenyl, 2,2',5,5'-tetrach...	290	C12H6Cl4	035693-99-3	99		
4	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99		
5	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
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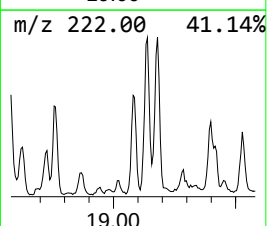
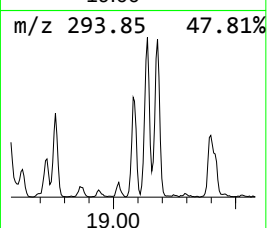
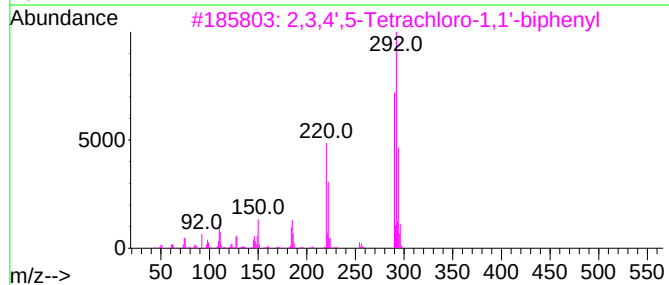
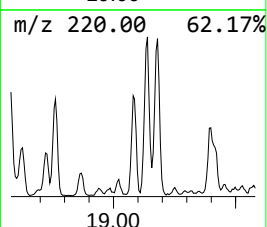
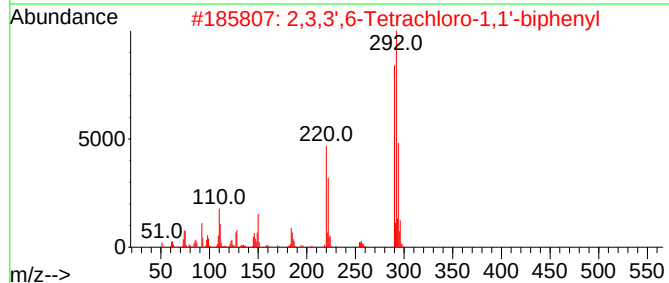
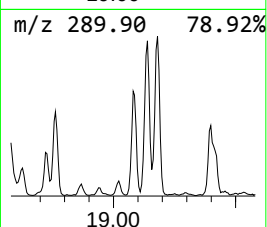
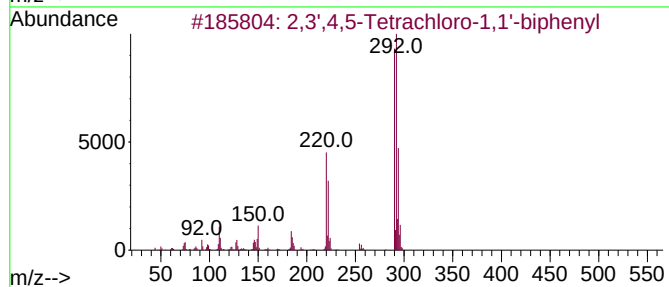
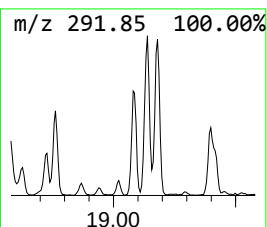
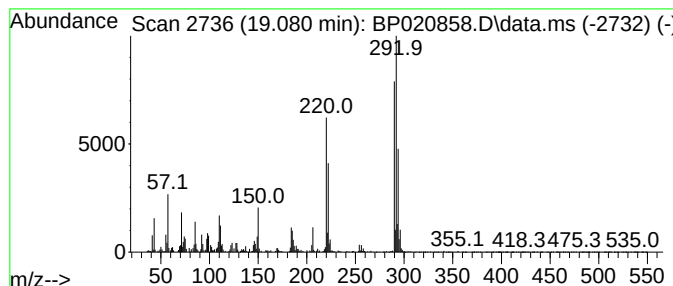
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.080	3.44 ng/ul	542271	Phenanthrene-d10	17.174

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	99	
2	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	99	
3	2,3,4',5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-34-7	99	
4	1,1'-Biphenyl, 3,3',4,4'-tetrach...	290	C12H6Cl4	032598-13-3	99	
5	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

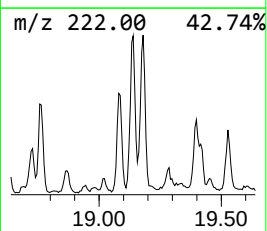
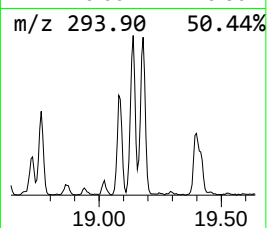
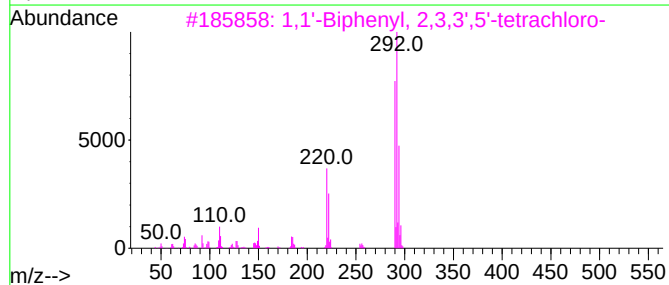
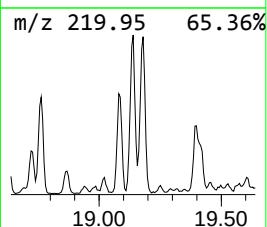
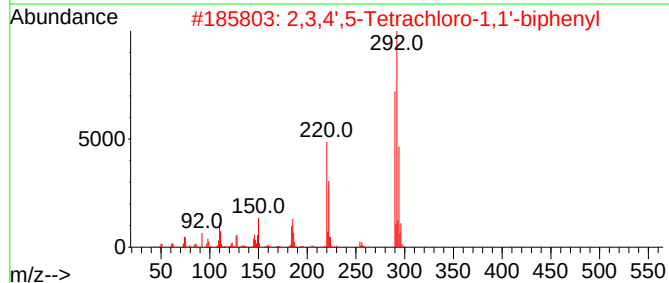
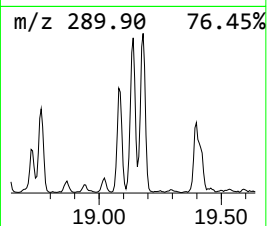
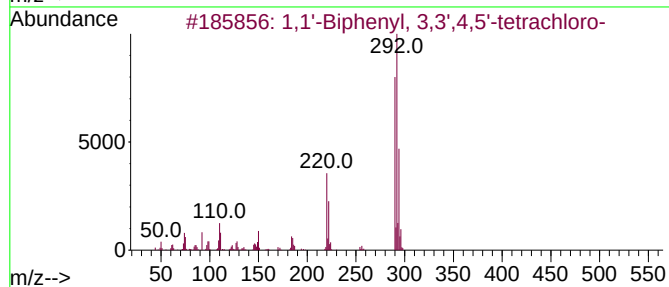
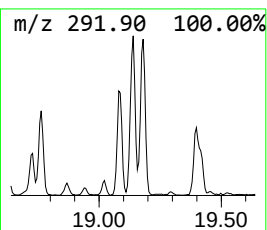
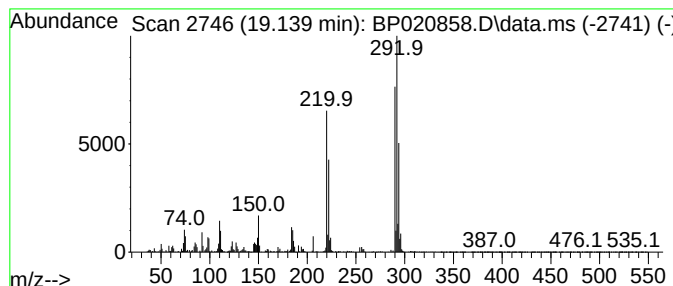
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,1'-Biphenyl, 3,3',4,5'-te... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.139	4.28 ng/ul	675077	Phenanthrene-d10	17.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	99
2	2,3,4',5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-34-7	99
3	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	99
4	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
5	Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
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Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
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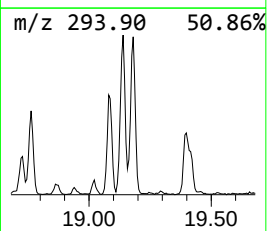
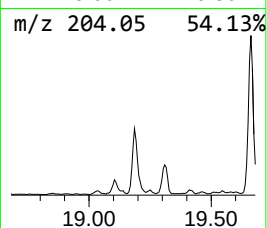
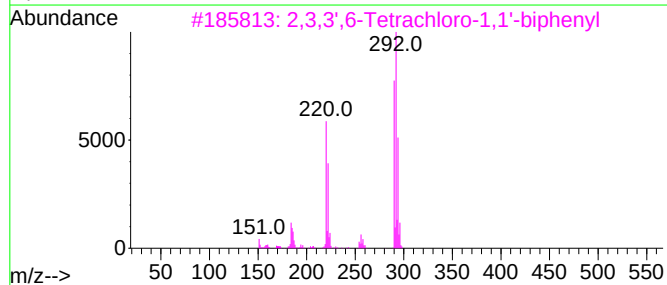
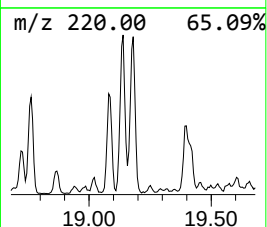
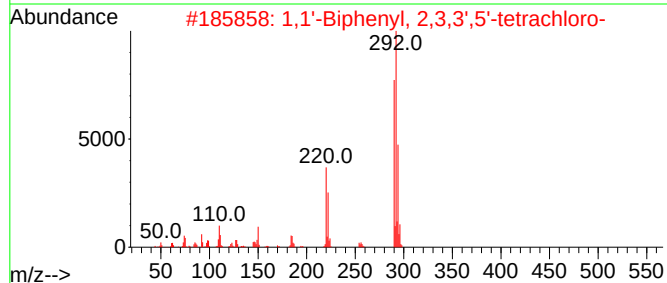
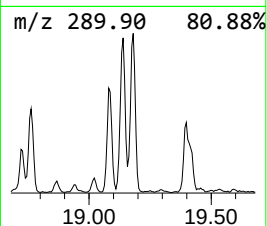
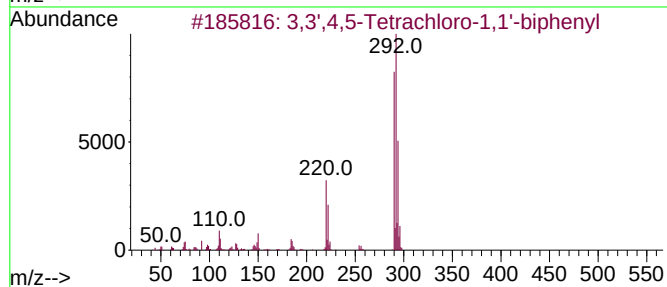
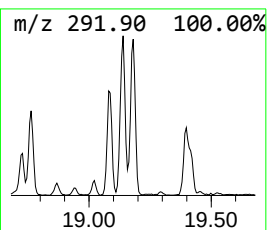
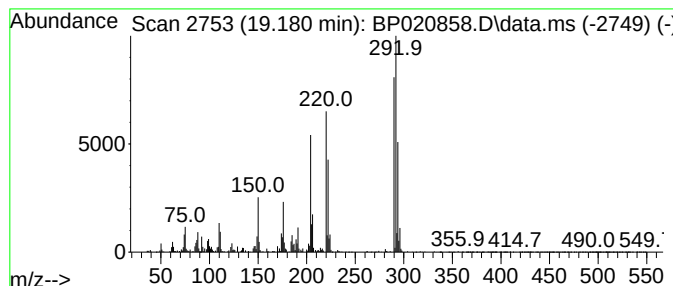
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 3,3',4,5-Tetrachloro-1,1'-b... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.180	5.66 ng/ul	892017	Phenanthrene-d10	17.174	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	95
2	1,1'-Biphenyl, 2,3,3',5'-tetrach...	290	C12H6Cl4	041464-49-7	94
3	2,3,3',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074472-33-6	94
4	1,1'-Biphenyl, 2,3',5,5'-tetrach...	290	C12H6Cl4	041464-42-0	93
5	1,1'-Biphenyl, 2,3',4',5-tetrach...	290	C12H6Cl4	032598-11-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Instrument :
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ClientSampleId :
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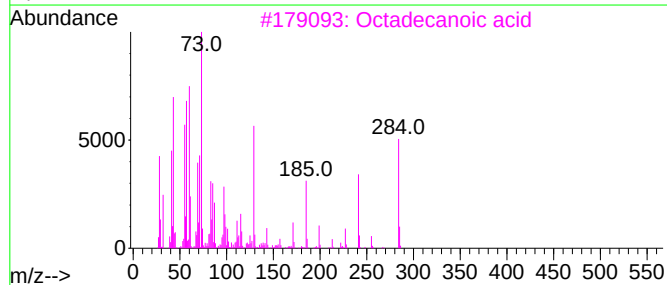
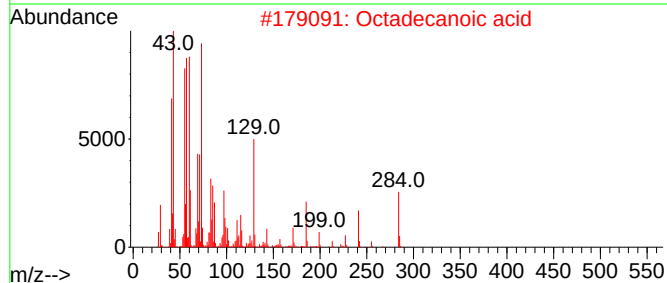
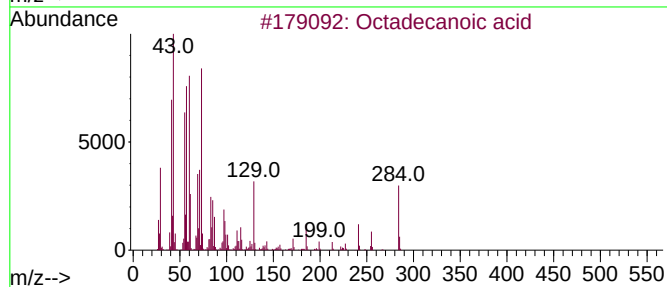
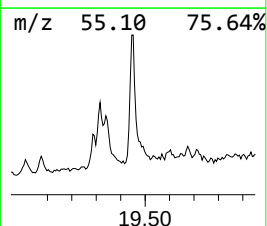
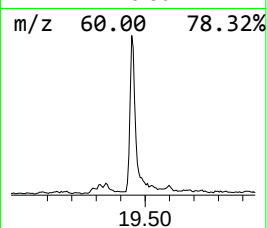
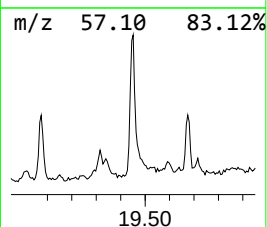
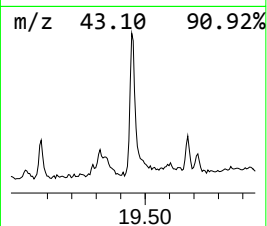
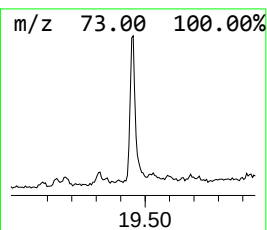
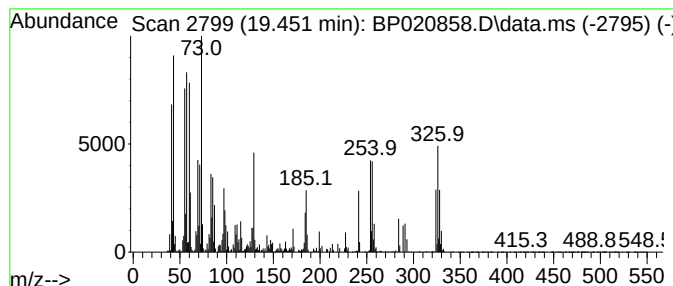
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Octadecanoic acid Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.451	3.93 ng/ul	1043610	Chrysene-d12	21.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	95
2			Octadecanoic acid	284	C18H36O2	000057-11-4	94
3			Octadecanoic acid	284	C18H36O2	000057-11-4	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	93
5			Octadecanoic acid	284	C18H36O2	000057-11-4	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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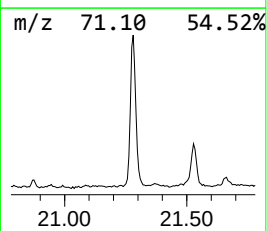
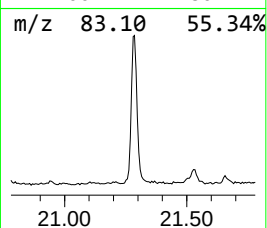
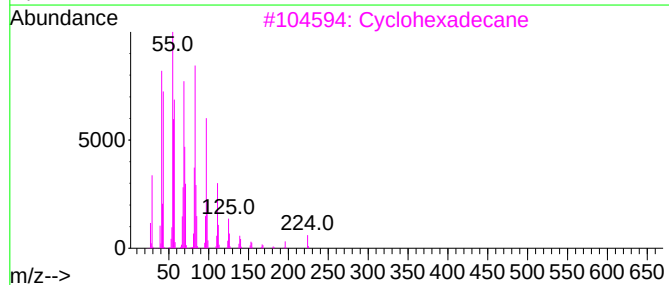
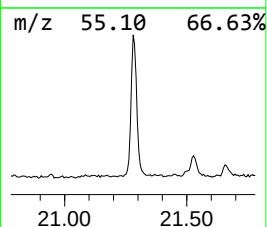
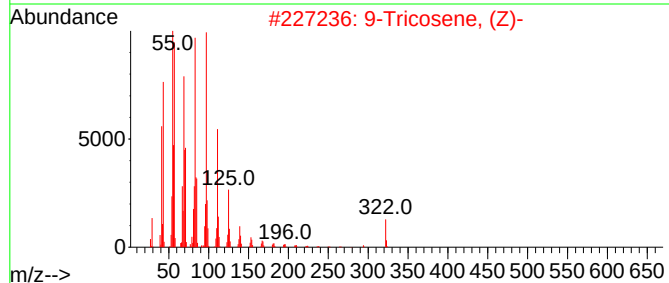
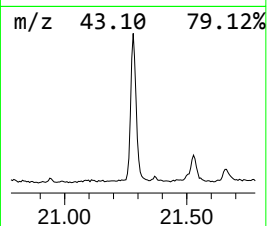
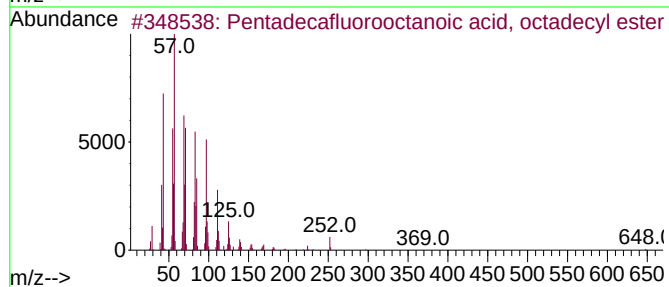
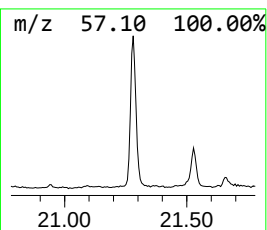
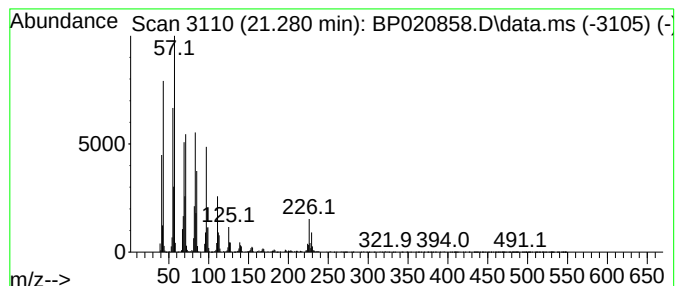
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Pentadecafluorooctanoic aci... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.280	10.69 ng/ul	2841110	Chrysene-d12	21.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2			9-Tricosene, (Z)-	322	C23H46	027519-02-4	92
3			Cyclohexadecane	224	C16H32	000295-65-8	91
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	90
5			17-Pentatriacontene	491	C35H70	006971-40-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Misc :
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Instrument :
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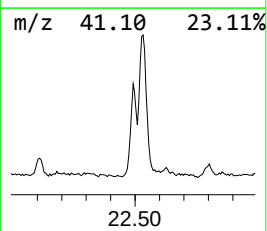
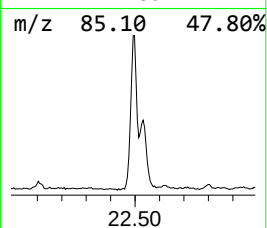
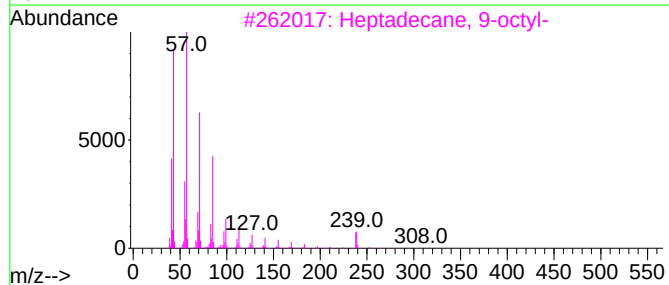
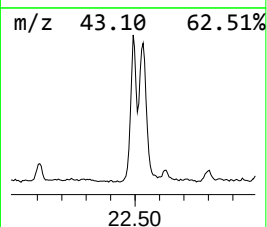
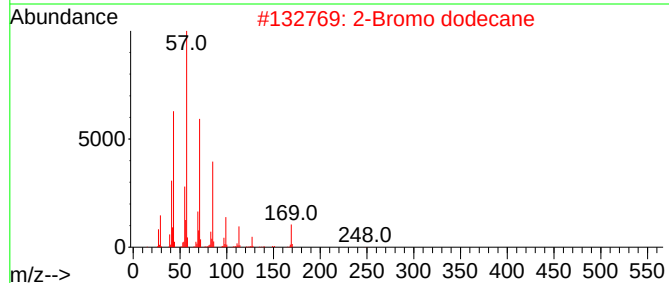
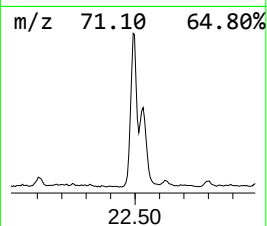
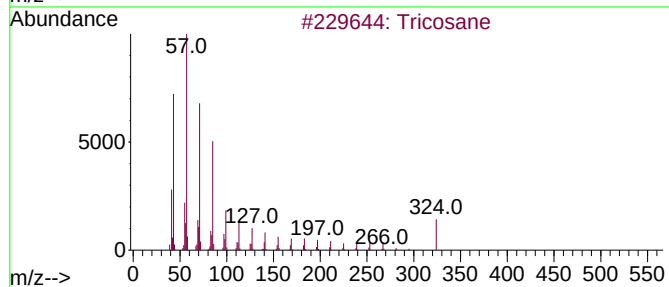
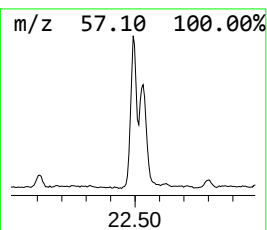
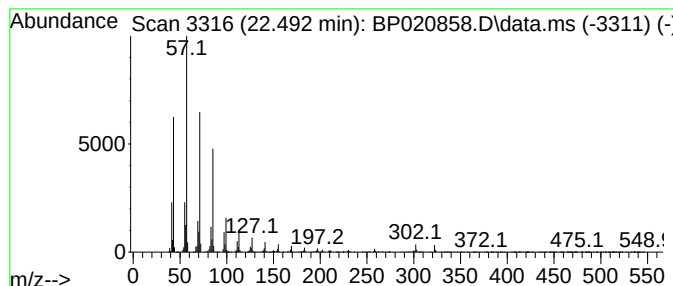
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.492	6.32 ng/ul	1680680	Chrysene-d12	21.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	94
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	91
3			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4			Heptacosane	380	C27H56	000593-49-7	90
5			Octacosane, 2-methyl-	408	C29H60	001560-98-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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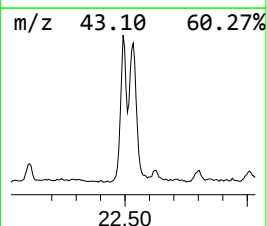
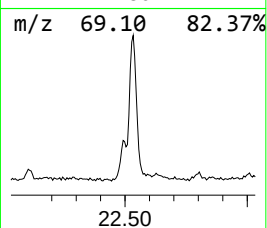
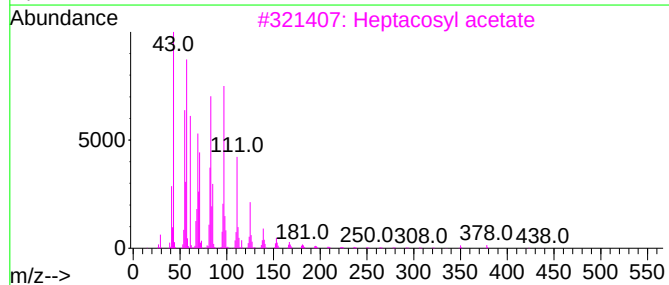
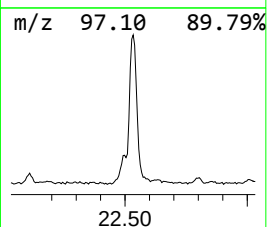
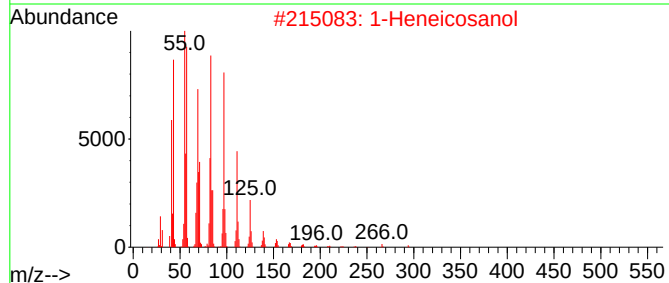
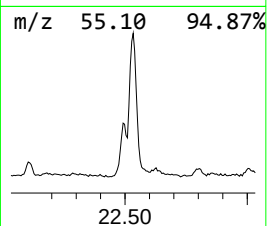
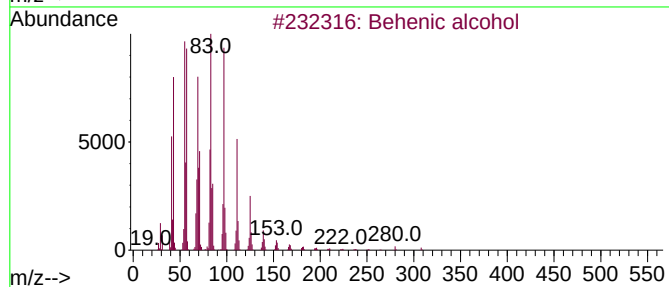
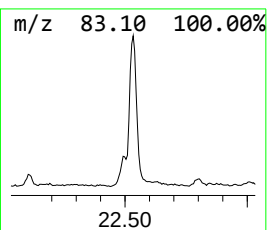
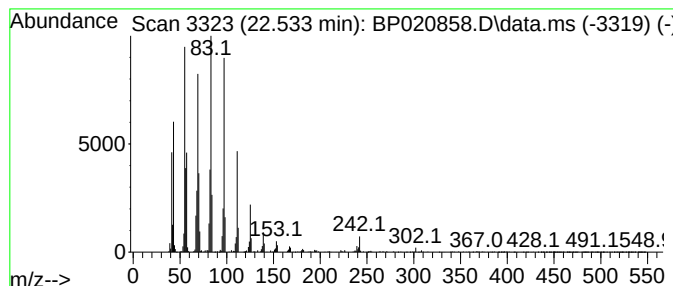
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Behenic alcohol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.533	11.26 ng/ul	2994650	Chrysene-d12	21.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	94
3			Heptacosyl acetate	438	C29H58O2	1000351-78-2	74
4			Cyclopentadecane	210	C15H30	000295-48-7	70
5			1-Octadecanol	270	C18H38O	000112-92-5	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
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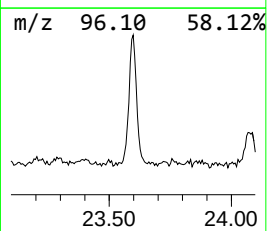
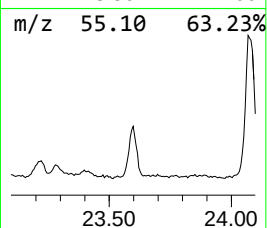
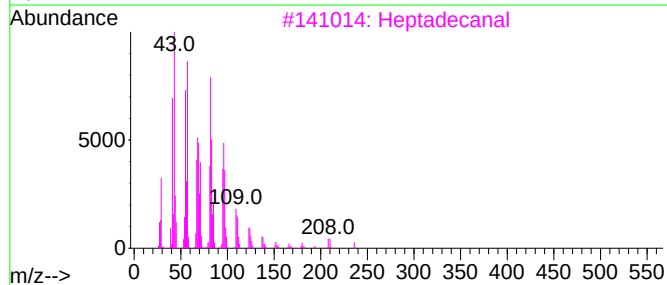
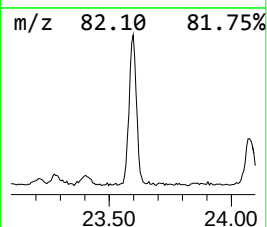
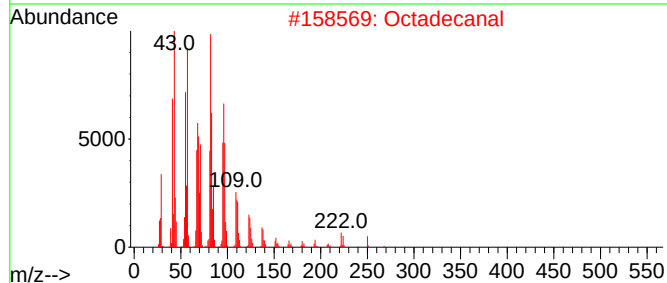
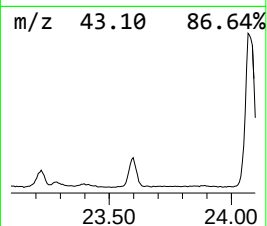
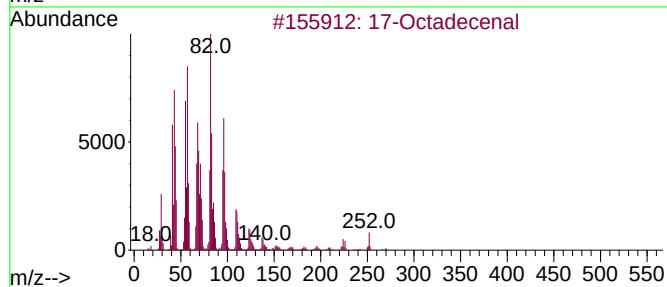
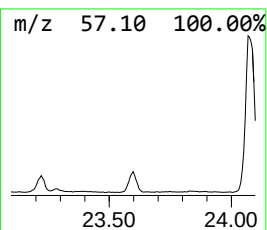
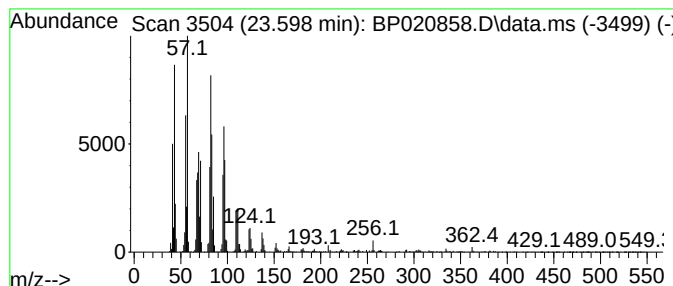
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 17-Octadecenal Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.598	9.92 ng/ul	1474380	Perylene-d12	24.968

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		17-Octadecenal	266	C18H34O	056554-86-0	91
2		Octadecanal	268	C18H36O	000638-66-4	91
3		Heptadecanal	254	C17H34O	1000376-70-0	91
4		16-Octadecenal	266	C18H34O	056554-87-1	91
5		Hexacosanal	380	C26H52O	026627-85-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

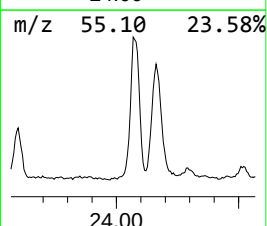
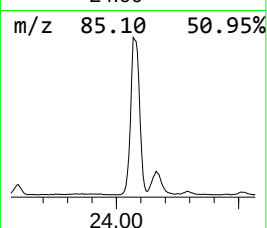
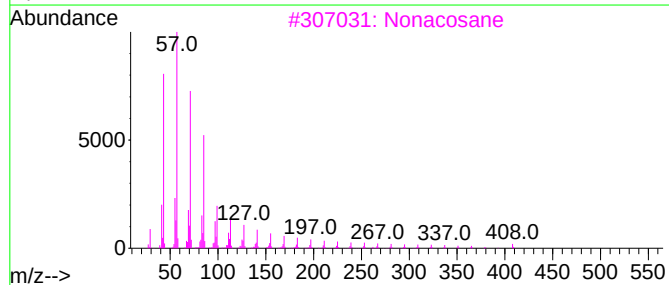
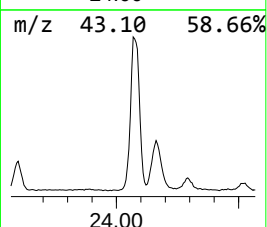
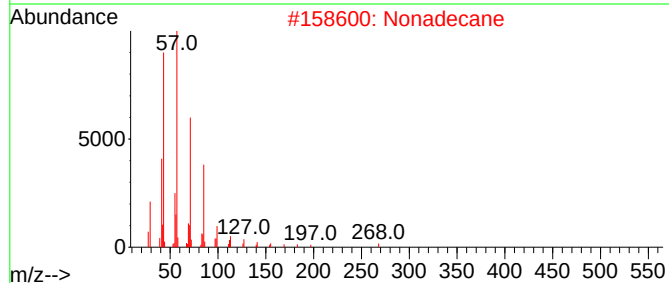
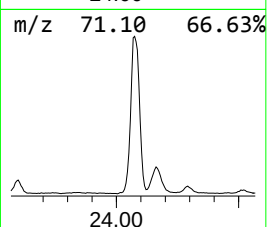
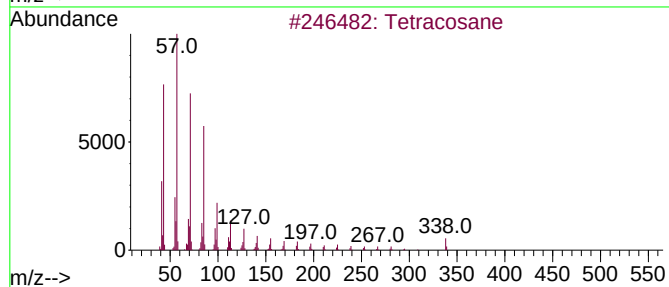
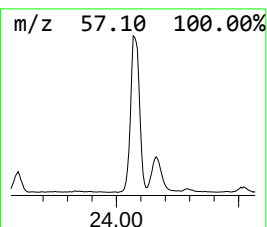
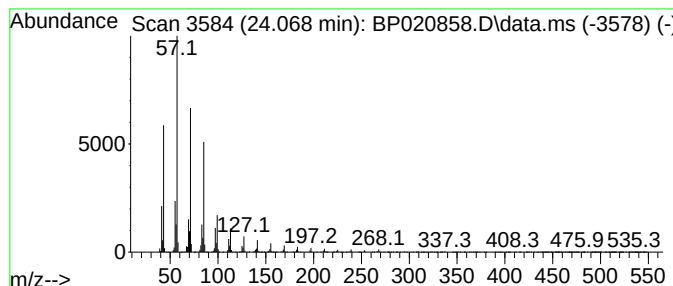
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.068	43.37 ng/ul	6448340	Perylene-d12	24.968	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Nonadecane	268	C19H40	000629-92-5	95
3	Nonacosane	408	C29H60	000630-03-5	94
4	Heptadecane, 8,8-dipentyl-	380	C27H56	055282-28-5	94
5	2-Methylpentacosane	366	C26H54	000629-87-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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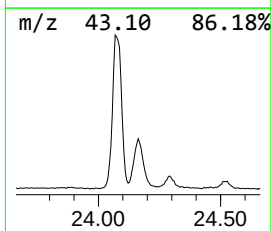
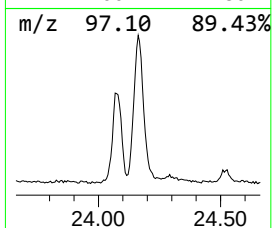
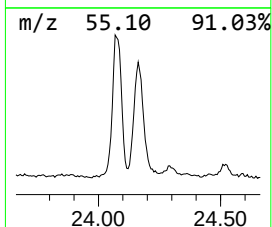
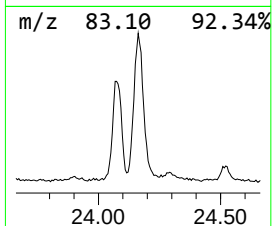
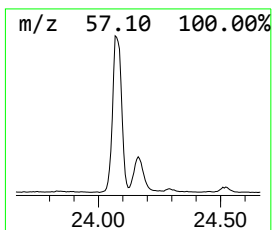
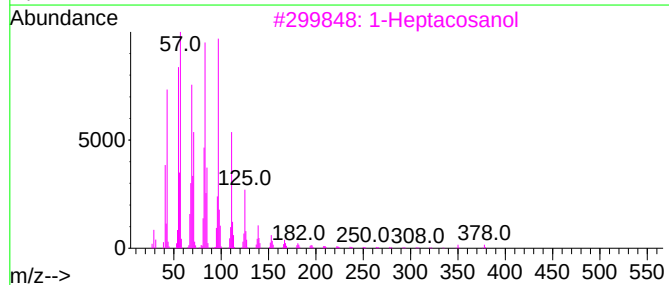
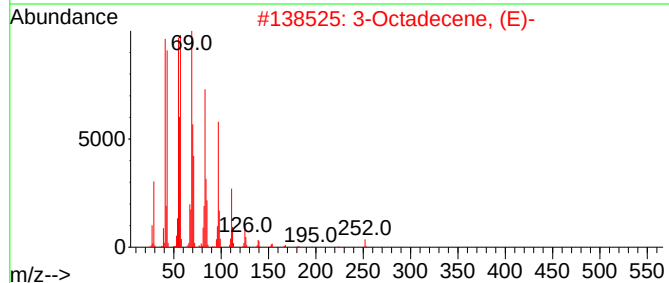
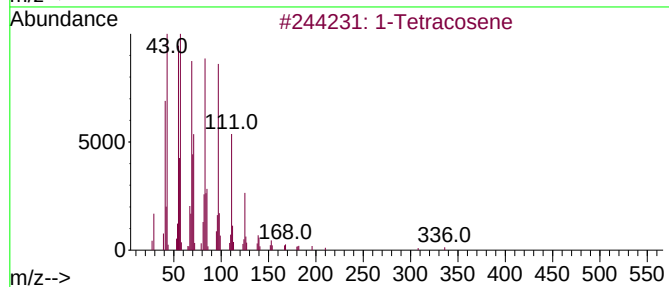
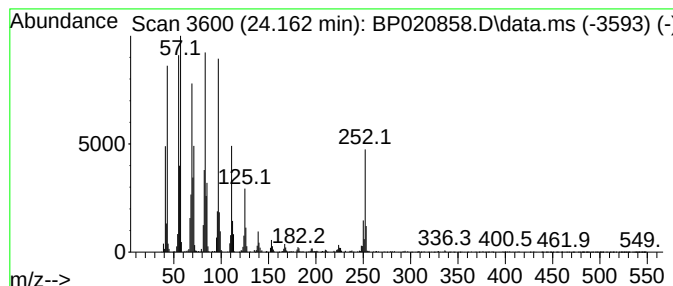
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.162	23.65 ng/ul	3516740	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tetracosene			336	C24H48	010192-32-2	95
2	3-Octadecene, (E)-			252	C18H36	007206-19-1	94
3	1-Heptacosanol			396	C27H56O	002004-39-9	93
4	1-Heneicosanol			312	C21H44O	015594-90-8	93
5	Nonadecyl trifluoroacetate			380	C21H39F3O2	1000351-76-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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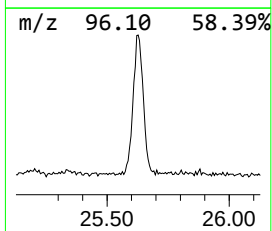
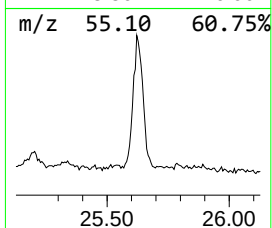
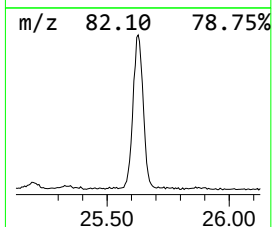
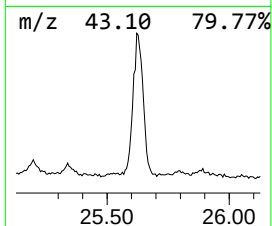
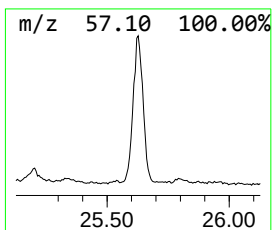
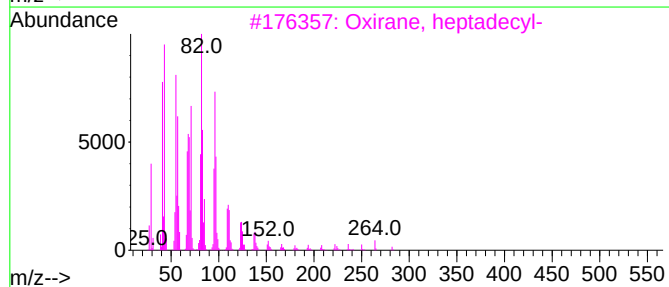
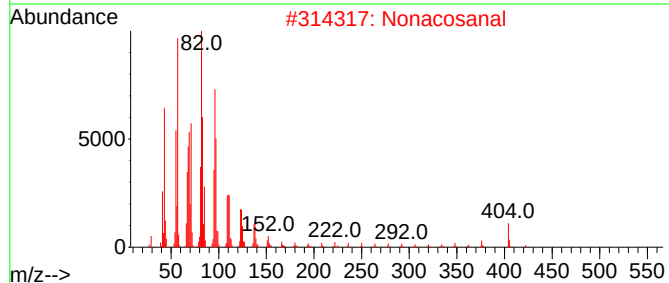
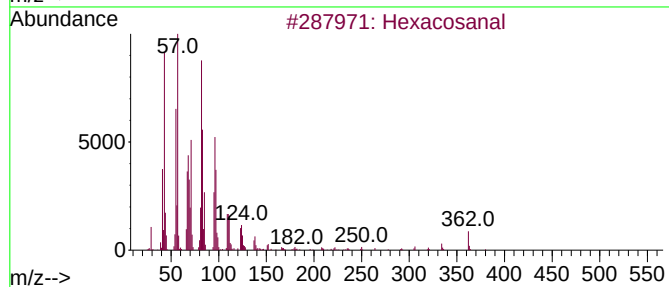
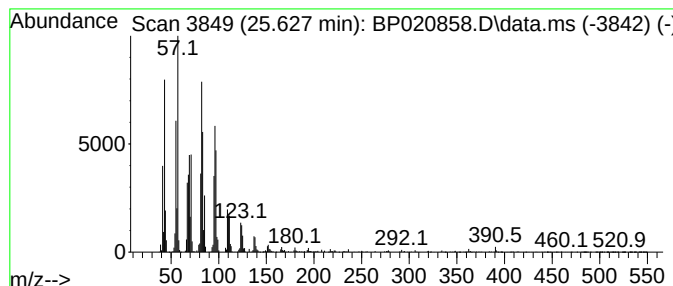
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 Hexacosanal Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.627	21.53 ng/ul	3201880	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosanal	380	C26H52O	026627-85-0	94
2			Nonacosanal	422	C29H58O	072934-04-4	93
3			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91
4			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
5			Tetracosanal	352	C24H48O	057866-08-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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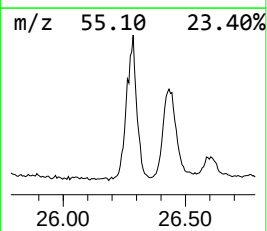
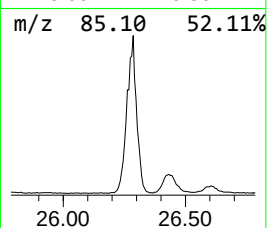
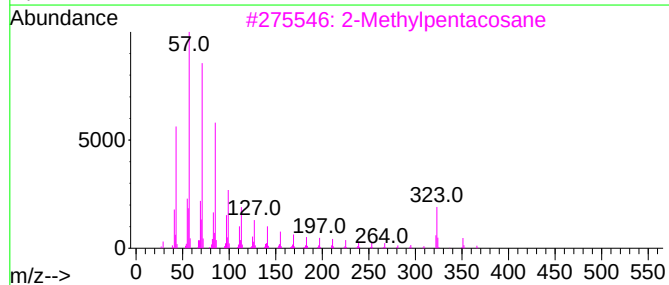
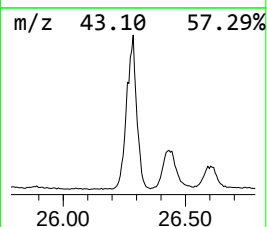
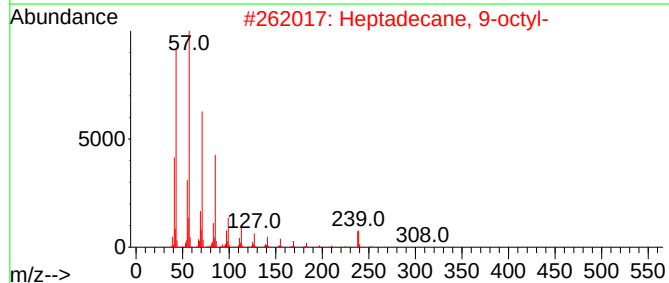
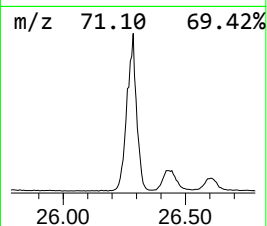
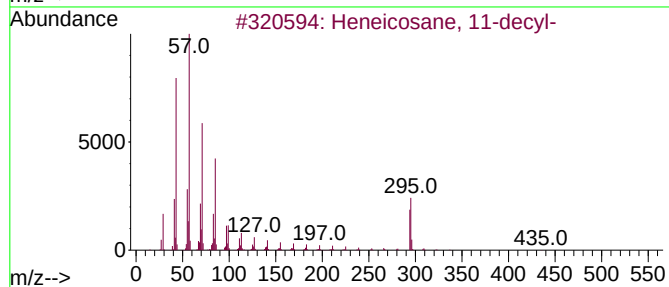
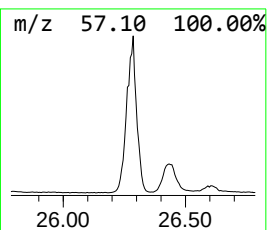
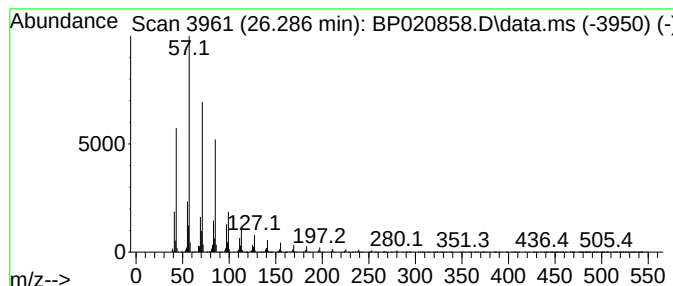
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.286	43.78 ng/ul	6509270	Perylene-d12	24.968

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	97
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3		2-Methylpentacosane	366	C26H54	000629-87-8	94
4		Heptacosane	380	C27H56	000593-49-7	91
5		Hexacosane	366	C26H54	000630-01-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Sample : P3035-01
Misc :
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Instrument :
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ClientSampleId :
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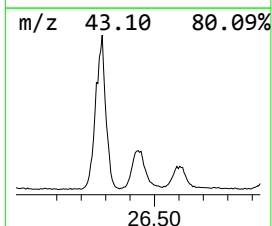
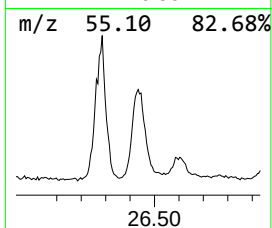
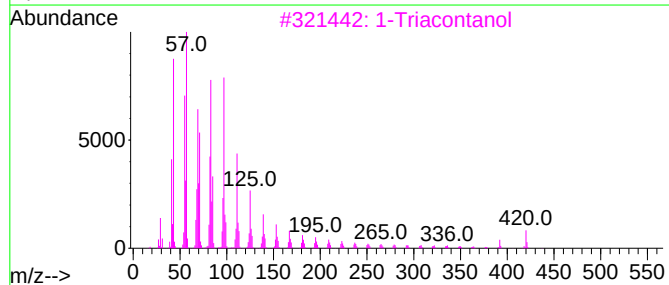
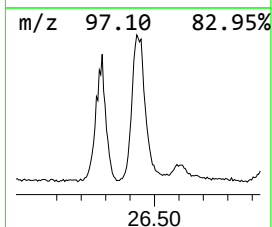
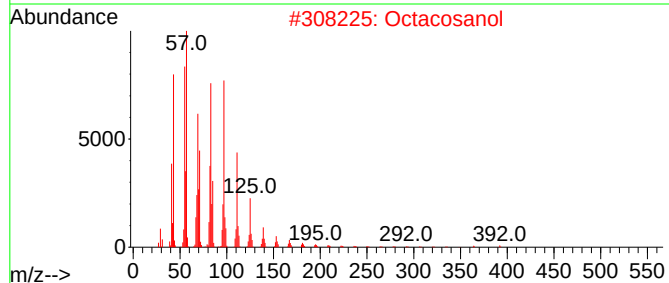
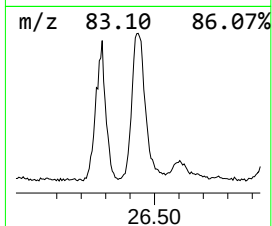
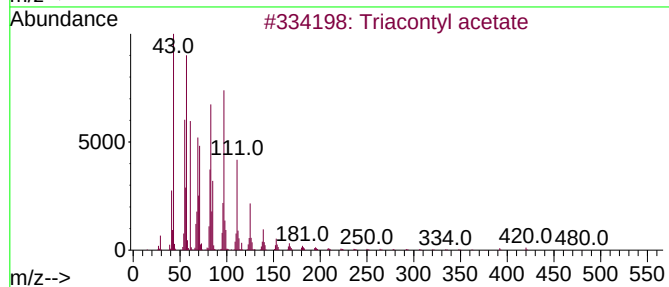
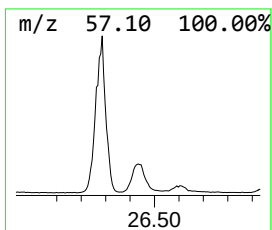
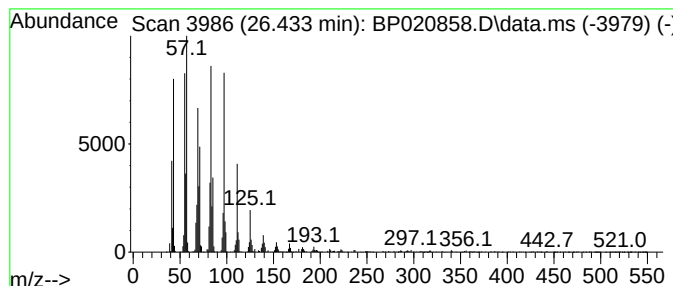
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Triacontyl acetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.433	8.84 ng/ul	1313830	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triacontyl acetate	480	C32H64O2	041755-58-2	97
2			Octacosanol	410	C28H58O	000557-61-9	95
3			1-Triacontanol	438	C30H62O	000593-50-0	94
4			1-Hexacosene	364	C26H52	018835-33-1	94
5			2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
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Sample : P3035-01
Misc :
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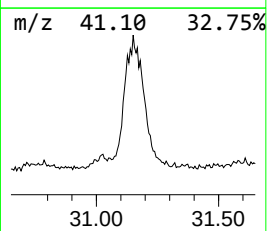
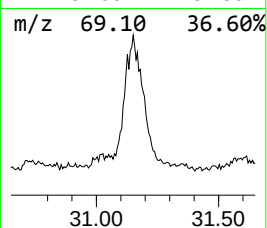
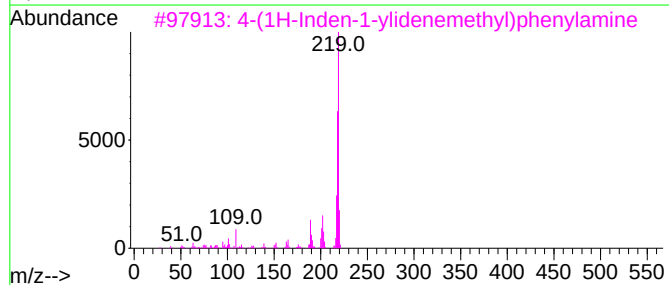
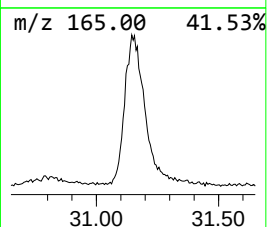
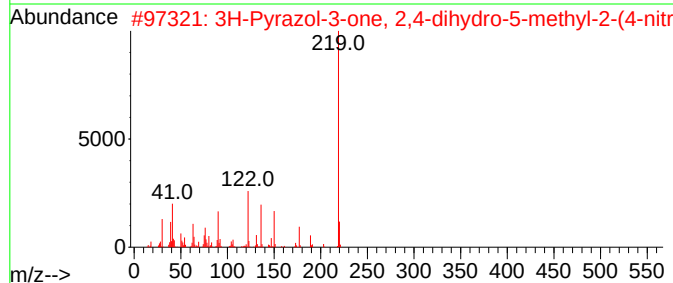
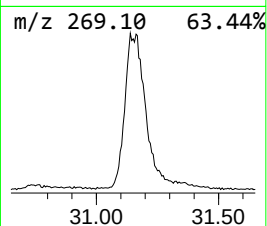
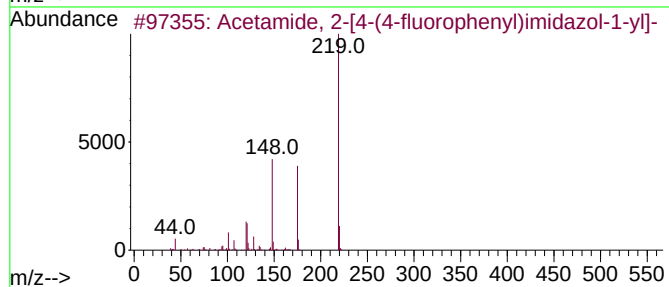
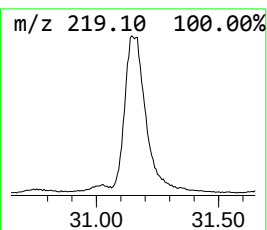
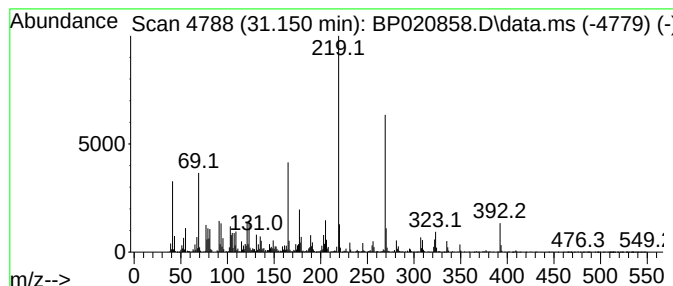
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
31.150	13.28 ng/ul	1974260	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, 2-[4-(4-fluorophenyl)...	219	C11H10FN3O	1000303-92-2	38
2			3H-Pyrazol-3-one, 2,4-dihydro-5-...	219	C10H9N3O3	006402-09-1	35
3			4-(1H-Inden-1-ylidenemethyl)phen...	219	C16H13N	000487-61-6	30
4			8-(2-Hydroxy-1-methoxy-3-methylb...	304	C17H20O5	1000493-42-6	27
5			3,4-Dihydrocoumarin, 4,4-dimethy...	219	C13H17NO2	029423-77-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070824\
Data File : BP020858.D
Acq On : 09 Jul 2024 10:25
Operator : MA/JU
Sample : P3035-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

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					#	RT	Resp	Conc
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unknown-02	4.881	2.4	ng/ul	143504	1	7.722	1204720	20.0
1,1'-Biphenyl, ...	17.804	7.1	ng/ul	1122010	4	17.174	3152430	20.0
n-Hexadecanoic ...	18.116	13.2	ng/ul	2080950	4	17.174	3152430	20.0
1,1'-Biphenyl, ...	18.286	6.5	ng/ul	1019290	4	17.174	3152430	20.0
1,1'-Biphenyl, ...	18.345	3.3	ng/ul	516864	4	17.174	3152430	20.0
1,1'-Biphenyl, ...	18.392	3.1	ng/ul	485238	4	17.174	3152430	20.0
1,1'-Biphenyl, ...	18.575	2.5	ng/ul	393510	4	17.174	3152430	20.0
2,3',4,5-Tetrac...	19.080	3.4	ng/ul	542271	4	17.174	3152430	20.0
1,1'-Biphenyl, ...	19.139	4.3	ng/ul	675077	4	17.174	3152430	20.0
3,3',4,5-Tetrac...	19.180	5.7	ng/ul	892017	4	17.174	3152430	20.0
Octadecanoic acid	19.451	3.9	ng/ul	1043610	5	21.621	5317330	20.0
Pentadecafluoro...	21.280	10.7	ng/ul	2841110	5	21.621	5317330	20.0
(DEL) Alkane: S...	22.492	6.3	ng/ul	1680680	5	21.621	5317330	20.0
Behenic alcohol	22.533	11.3	ng/ul	2994650	5	21.621	5317330	20.0
17-Octadecenal	23.598	9.9	ng/ul	1474380	6	24.968	2973770	20.0
(DEL) Alkane: S...	24.068	43.4	ng/ul	6448340	6	24.968	2973770	20.0
(DEL) Alkane: S...	24.162	23.6	ng/ul	3516740	6	24.968	2973770	20.0
Hexacosanal	25.627	21.5	ng/ul	3201880	6	24.968	2973770	20.0
(DEL) Alkane: S...	26.286	43.8	ng/ul	6509270	6	24.968	2973770	20.0
Triacetyl acetate	26.433	8.8	ng/ul	1313830	6	24.968	2973770	20.0
unknown-03	31.150	13.3	ng/ul	1974260	6	24.968	2973770	20.0