

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
 Data File : BP020807.D
 Acq On : 06 Jul 2024 04:50
 Operator : MA/JU
 Sample : P2964-12
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CS6

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: BP020807.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	53	rVB	30397	43345	0.93%	0.066%
2	3.534	88	93	101	rBV	37633	61239	1.32%	0.093%
3	3.687	114	119	130	rBV3	49478	85087	1.83%	0.129%
4	3.775	130	134	140	rVB	52439	67680	1.46%	0.103%
5	3.987	166	170	174	rVB2	19774	26265	0.57%	0.040%
6	4.534	259	263	267	rVV	32061	44243	0.95%	0.067%
7	4.581	267	271	276	rVB	31839	42818	0.92%	0.065%
8	4.881	316	322	330	rBV	157787	239886	5.16%	0.364%
9	5.846	477	486	493	rBV2	23632	45235	0.97%	0.069%
10	6.916	658	668	682	rBV	567328	973483	20.95%	1.476%
11	7.069	687	694	707	rBV	429981	711904	15.32%	1.079%
12	7.269	720	728	737	rBV	622267	984221	21.18%	1.492%
13	7.722	794	805	813	rBV	687533	1155379	24.87%	1.751%
14	7.799	813	818	827	rVB3	21996	44353	0.95%	0.067%
15	8.216	883	889	894	rBV3	12365	22494	0.48%	0.034%
16	8.434	919	926	934	rBV	603378	1020634	21.97%	1.547%
17	8.493	934	936	940	rVV4	14226	21260	0.46%	0.032%
18	8.546	940	945	954	rVB2	27149	59611	1.28%	0.090%
19	8.881	994	1002	1016	rBV	547297	948744	20.42%	1.438%
20	9.252	1059	1065	1073	rBV3	20668	38723	0.83%	0.059%
21	9.434	1088	1096	1102	rBV	77688	121814	2.62%	0.185%
22	9.593	1114	1123	1137	rBV	456506	833017	17.93%	1.263%
23	10.134	1208	1215	1224	rBV	715446	1255538	27.02%	1.903%
24	10.499	1265	1277	1282	rBV	1007661	1693864	36.46%	2.568%
25	10.640	1294	1301	1318	rBV	147702	340169	7.32%	0.516%
26	11.040	1362	1369	1377	rBV2	23192	45863	0.99%	0.070%
27	11.240	1396	1403	1410	rBV2	77406	161284	3.47%	0.244%
28	11.904	1512	1516	1522	rVB2	18530	31927	0.69%	0.048%
29	12.098	1541	1549	1553	rVV2	12950	26985	0.58%	0.041%
30	12.169	1557	1561	1569	rVB	14241	26028	0.56%	0.039%
31	12.369	1591	1595	1605	rVB2	10921	24751	0.53%	0.038%
32	13.163	1724	1730	1739	rVV2	27771	47283	1.02%	0.072%
33	13.287	1747	1751	1757	rBV6	9630	21029	0.45%	0.032%
34	13.345	1757	1761	1778	rVB	47552	97973	2.11%	0.149%
35	13.522	1784	1791	1801	rBV2	40315	123522	2.66%	0.187%
36	13.669	1810	1816	1821	rBV3	25925	44739	0.96%	0.068%

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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

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	13.763	1823	1832	1837	rBV	1430720	1894027	40.76%	2.871%
38	14.045	1869	1880	1893	rBV	1579204	2426881	52.23%	3.679%
39	14.181	1898	1903	1908	rVV5	28291	53616	1.15%	0.081%
40	14.240	1908	1913	1918	rVB3	18305	28880	0.62%	0.044%
41	14.357	1923	1933	1940	rBV	1456846	2239698	48.20%	3.395%
42	14.422	1940	1944	1951	rVB	68722	96970	2.09%	0.147%
43	14.598	1963	1974	1994	rBV3	468446	1094163	23.55%	1.659%
44	14.745	1994	1999	2001	rBV	40972	74987	1.61%	0.114%
45	14.822	2009	2012	2020	rVB2	49473	71741	1.54%	0.109%
46	15.039	2045	2049	2054	rVB5	13185	20099	0.43%	0.030%
47	15.163	2062	2070	2074	rBV6	17746	42471	0.91%	0.064%
48	15.369	2098	2105	2111	rVV	1779958	2824137	60.78%	4.281%
49	15.422	2111	2114	2122	rVB	50972	83922	1.81%	0.127%
50	15.504	2122	2128	2139	rBV	359986	660469	14.22%	1.001%
51	15.634	2146	2150	2154	rVB3	39406	59876	1.29%	0.091%
52	15.722	2161	2165	2172	rBV3	26293	41055	0.88%	0.062%
53	15.881	2188	2192	2199	rVB2	20848	38877	0.84%	0.059%
54	15.951	2199	2204	2210	rBV8	35435	73381	1.58%	0.111%
55	16.110	2225	2231	2238	rBV6	36689	89216	1.92%	0.135%
56	16.375	2273	2276	2280	rVB3	24980	33066	0.71%	0.050%
57	16.504	2294	2298	2303	rBV4	17619	30477	0.66%	0.046%
58	16.557	2303	2307	2313	rBV7	36715	80234	1.73%	0.122%
59	16.686	2327	2329	2333	rVB5	27934	29772	0.64%	0.045%
60	16.816	2347	2351	2358	rVB2	55418	87092	1.87%	0.132%
61	16.892	2360	2364	2366	rVV4	22481	38376	0.83%	0.058%
62	16.928	2366	2370	2375	rVV4	82165	139785	3.01%	0.212%
63	16.981	2375	2379	2383	rVV	60827	91416	1.97%	0.139%
64	17.039	2385	2389	2391	rVV2	80119	110476	2.38%	0.167%
65	17.069	2391	2394	2400	rVB3	63335	116070	2.50%	0.176%
66	17.169	2405	2411	2415	rBV	1760063	2676694	57.61%	4.057%
67	17.210	2415	2418	2423	rVB	968443	1281829	27.59%	1.943%
68	17.269	2423	2428	2433	rBV2	1769574	2929644	63.05%	4.441%
69	17.345	2438	2441	2449	rVB3	78895	143701	3.09%	0.218%
70	17.592	2479	2483	2487	rVB	83205	101802	2.19%	0.154%
71	17.792	2510	2517	2522	rVB2	274233	461687	9.94%	0.700%
72	17.904	2533	2536	2542	rVB6	65545	115790	2.49%	0.176%
73	18.069	2556	2564	2567	rBV2	169838	435846	9.38%	0.661%
74	18.110	2567	2571	2579	rVB2	1073369	1641912	35.34%	2.489%
75	18.275	2594	2599	2604	rBV2	384579	648117	13.95%	0.982%
76	18.333	2604	2609	2613	rVB2	130149	266609	5.74%	0.404%

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77	18.380	2614	2617	2624	rBV3	128491	201401	4.33%	0.305%
78	18.563	2645	2648	2651	rBV	154036	214499	4.62%	0.325%
79	18.645	2658	2662	2665	rVB	162971	205132	4.41%	0.311%
80	18.751	2677	2680	2685	rVB	105481	126410	2.72%	0.192%
81	19.075	2732	2735	2739	rVB3	121323	146905	3.16%	0.223%
82	19.127	2741	2744	2747	rBV	229124	262583	5.65%	0.398%
83	19.169	2748	2751	2759	rVB4	259002	431704	9.29%	0.654%
84	19.304	2768	2774	2781	rVB	2418964	3315111	71.35%	5.025%
85	19.445	2794	2798	2805	rBV3	564650	854948	18.40%	1.296%
86	19.651	2828	2833	2836	rBV	2534561	3869537	83.28%	5.866%
87	19.680	2836	2838	2847	rVB	1615777	2010711	43.28%	3.048%
88	19.927	2877	2880	2885	rBV2	116521	228219	4.91%	0.346%
89	20.204	2924	2927	2930	rVV	235673	280517	6.04%	0.425%
90	20.233	2930	2932	2939	rVB2	251946	373570	8.04%	0.566%
91	20.310	2941	2945	2948	rBV	193956	267131	5.75%	0.405%
92	21.286	3106	3111	3117	rVB3	1386182	1974166	42.49%	2.993%
93	21.527	3149	3152	3160	rVB	587974	849932	18.29%	1.288%
94	21.621	3161	3168	3173	rVV2	2132793	4646256	100.00%	7.043%
95	21.669	3173	3176	3188	rVB	1071601	1782613	38.37%	2.702%
96	22.510	3315	3319	3321	rBV2	526015	850506	18.31%	1.289%
97	23.910	3553	3557	3567	rBV2	601308	1496820	32.22%	2.269%
98	24.092	3582	3588	3595	rVB	921332	2238849	48.19%	3.394%
99	24.751	3694	3700	3707	rVB2	794254	1751474	37.70%	2.655%
100	24.968	3731	3737	3747	rVB	1057590	2683627	57.76%	4.068%

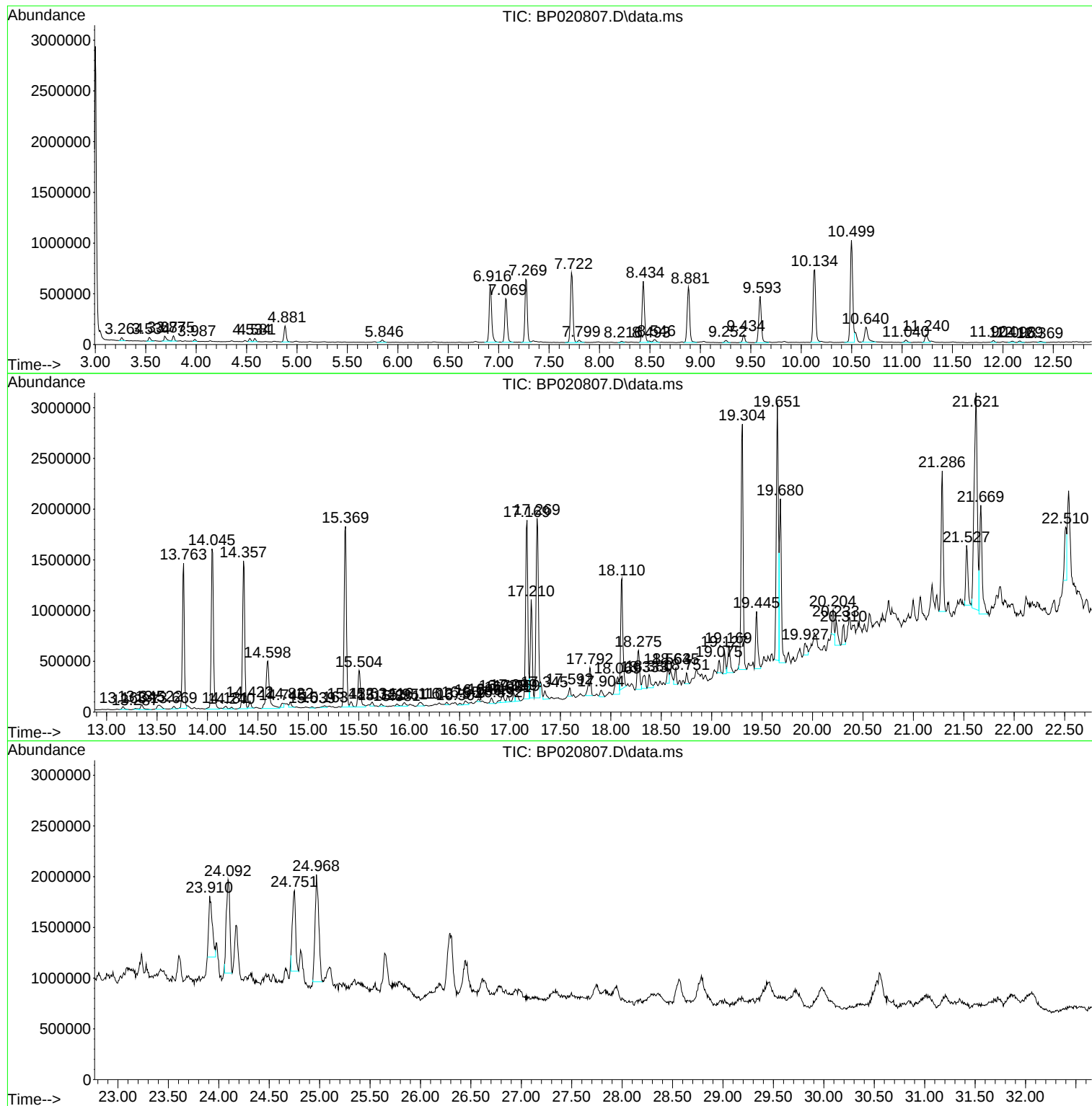
Sum of corrected areas: 65969872

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Instrument :
BNA_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



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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
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A4CS6

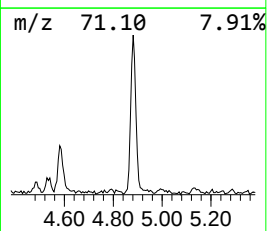
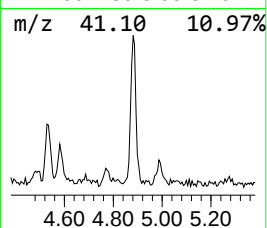
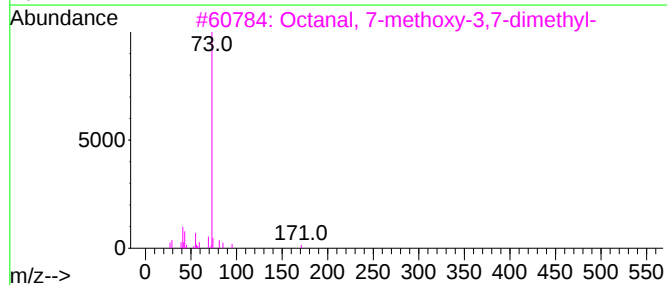
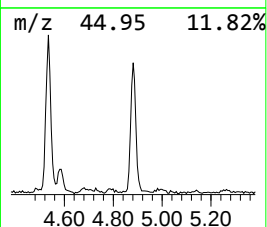
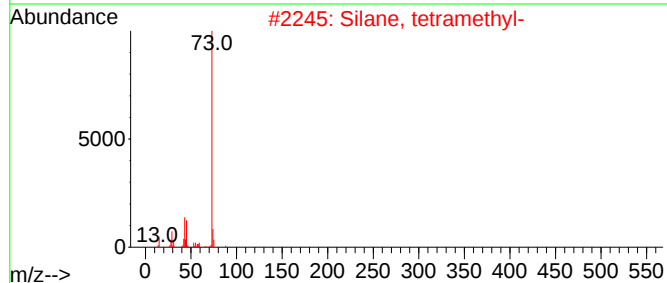
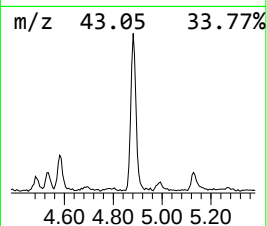
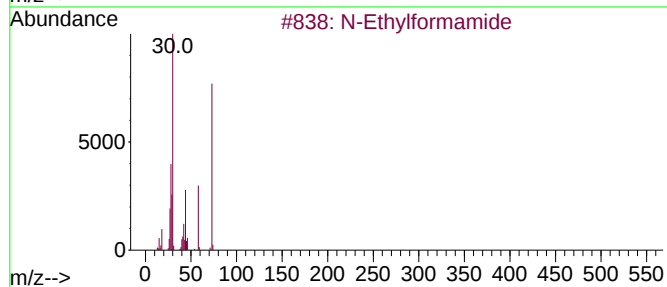
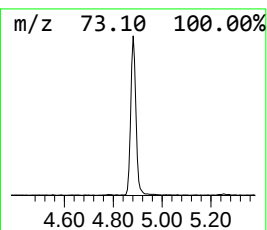
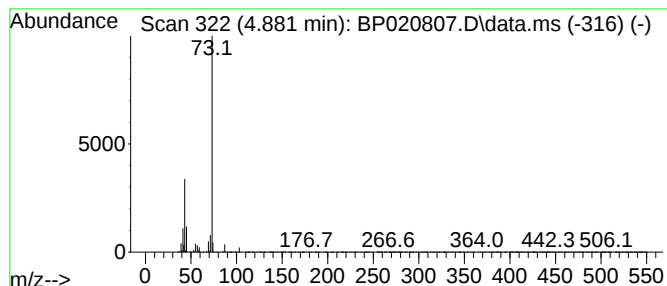
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.881	4.15 ng/ul	239886	1,4-Dichlorobenzene-d4	7.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Ethylformamide	73	C3H7NO	000627-45-2	9
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
3			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	9
4			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	9
5			Isopropyl 5,11-dihydroxy-3,7,11-...	314	C18H34O4	1000223-34-1	9



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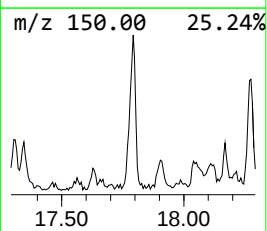
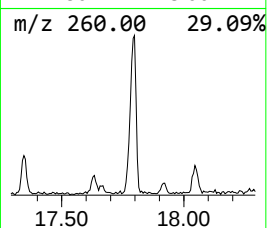
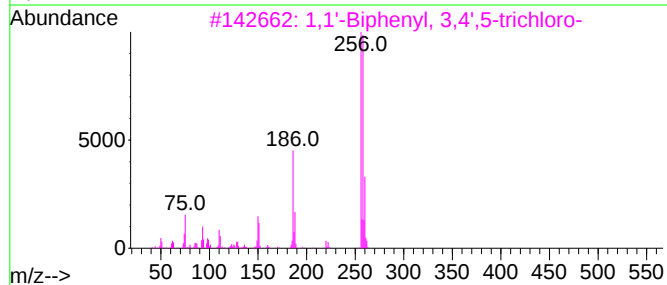
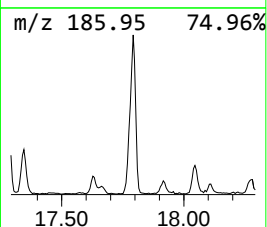
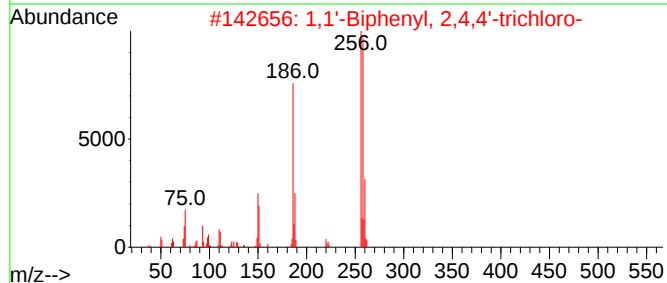
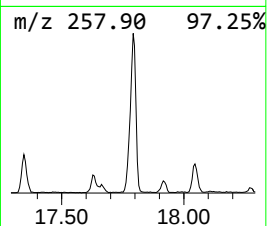
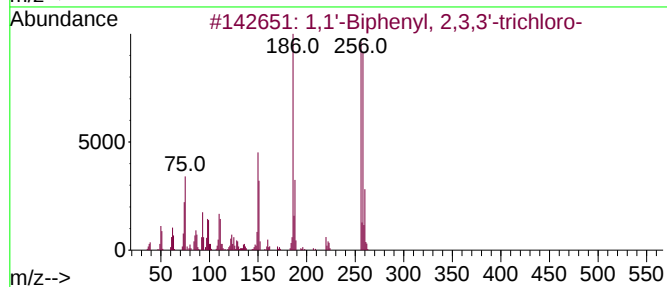
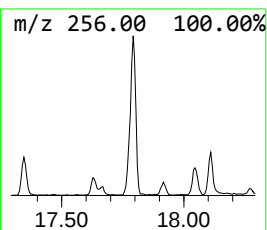
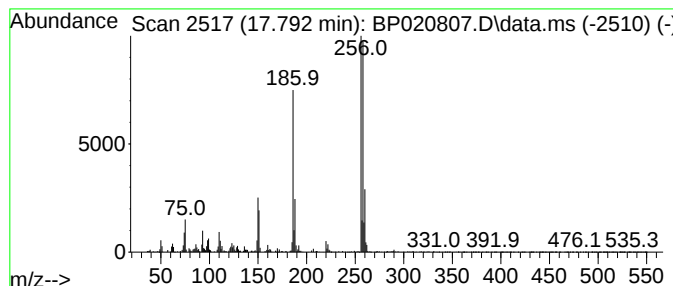
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,1'-Biphenyl, 2,3,3'-trich... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.792	3.45 ng/ul	461687	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3,3'-trichloro-	256	C12H7Cl3	038444-84-7	99		
2	1,1'-Biphenyl, 2,4,4'-trichloro-	256	C12H7Cl3	007012-37-5	97		
3	1,1'-Biphenyl, 3,4',5-trichloro-	256	C12H7Cl3	038444-88-1	96		
4	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	96		
5	1,1'-Biphenyl, 2,3,6-trichloro-	256	C12H7Cl3	055702-45-9	96		



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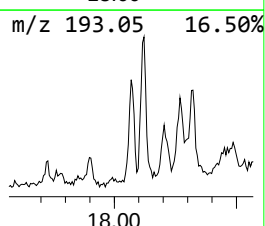
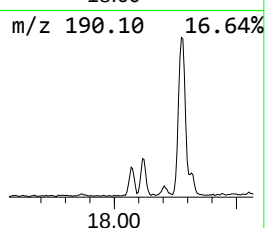
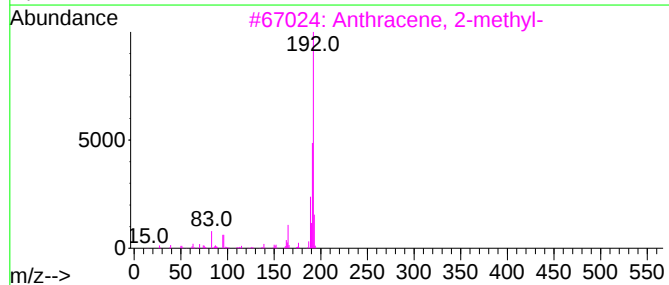
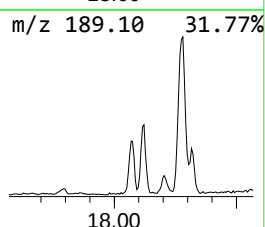
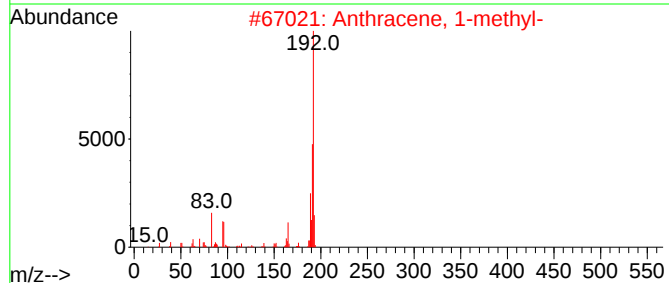
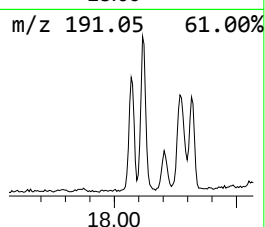
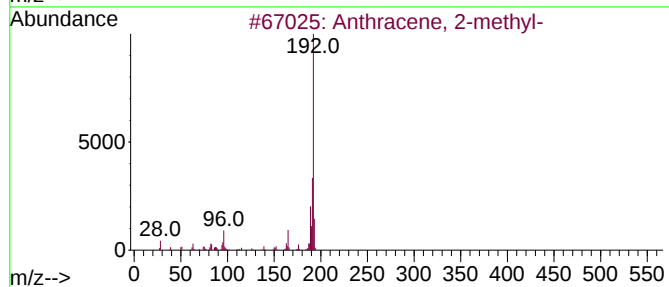
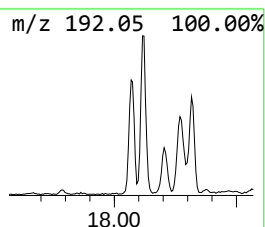
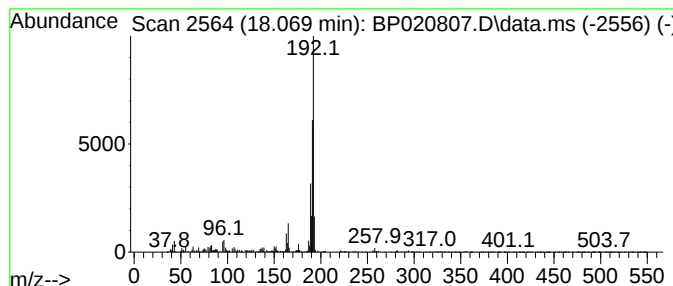
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Anthracene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.069	3.26 ng/ul	435846	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
2			Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93



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Acq On : 06 Jul 2024 04:50
Operator : MA/JU
Sample : P2964-12
Misc :
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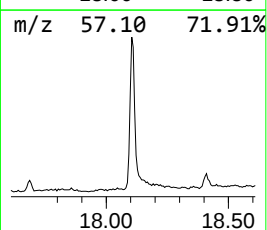
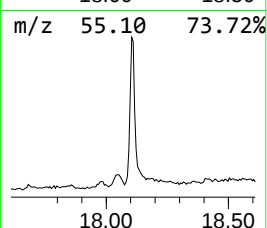
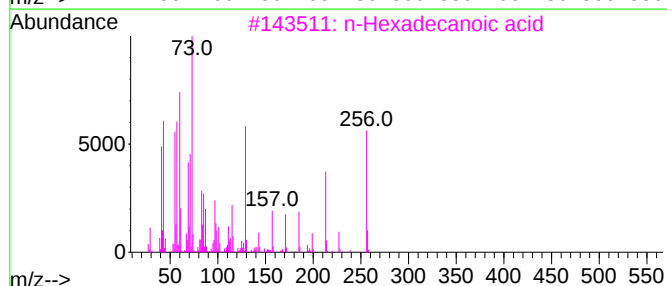
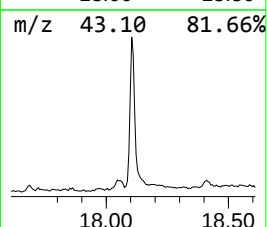
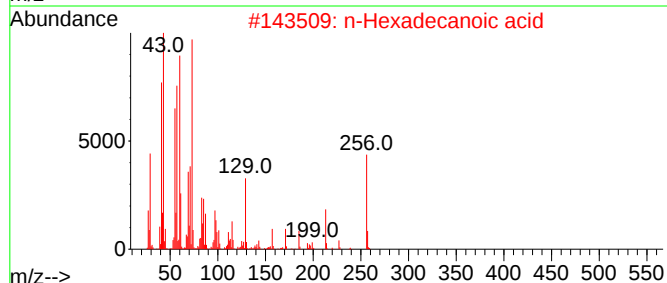
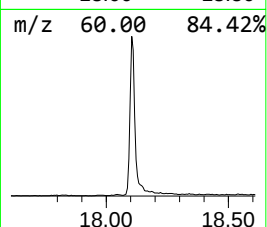
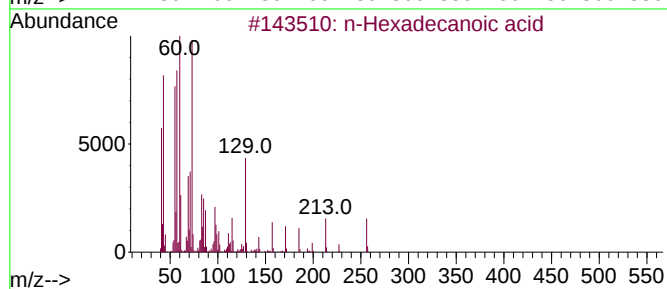
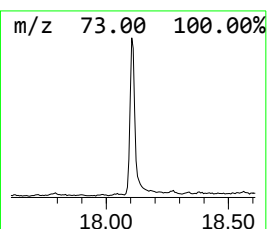
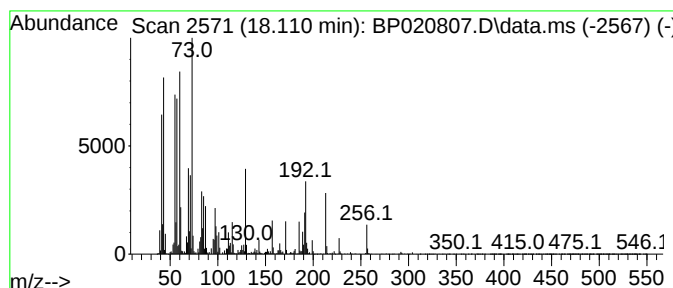
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.110	12.27 ng/ul	1641910	Phenanthrene-d10	17.169	
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	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	Tridecanoic acid	214	C13H26O2	000638-53-9	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020807.D
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Sample : P2964-12
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Instrument :
BNA_P
ClientSampleId :
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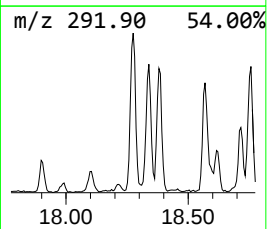
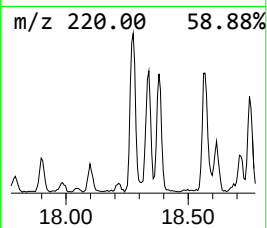
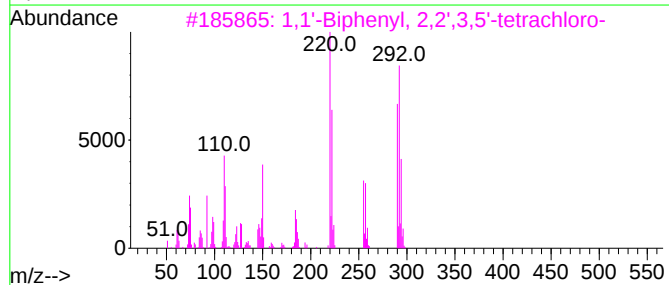
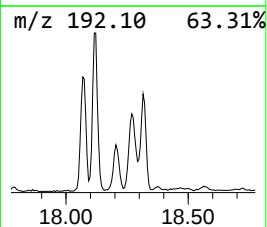
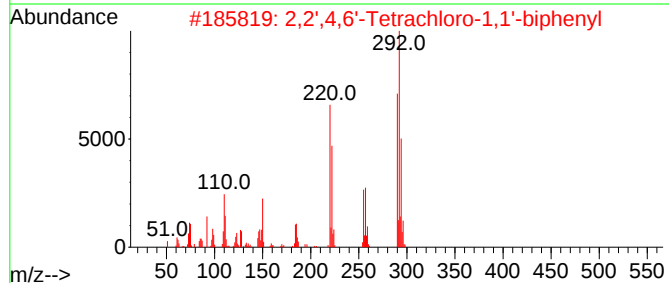
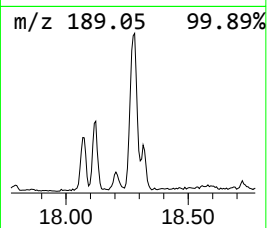
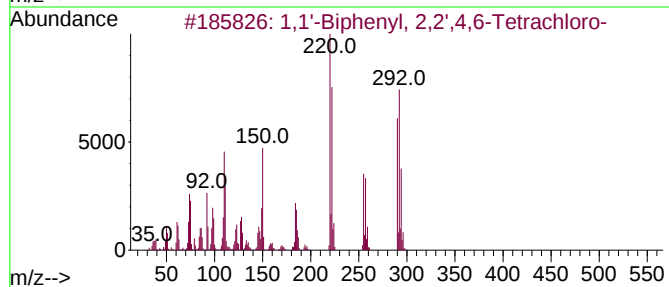
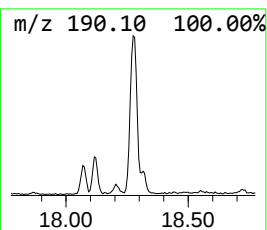
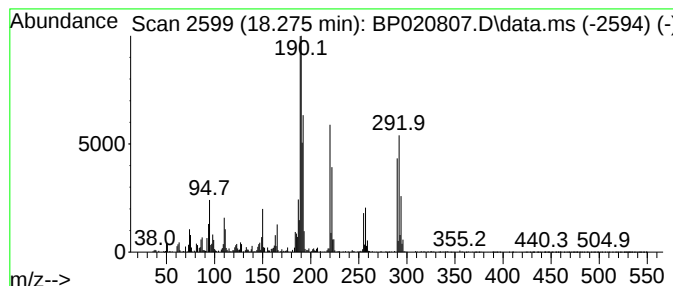
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.275	4.84 ng/ul	648117	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,6-Tetrachl...	290	C12H6Cl4	062796-65-0	96		
2	2,2',4,6'-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	068194-04-7	96		
3	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	96		
4	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	96		
5	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	95		



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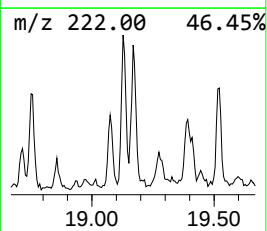
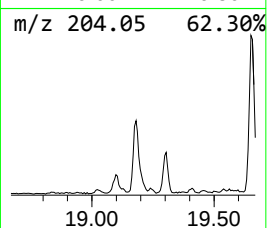
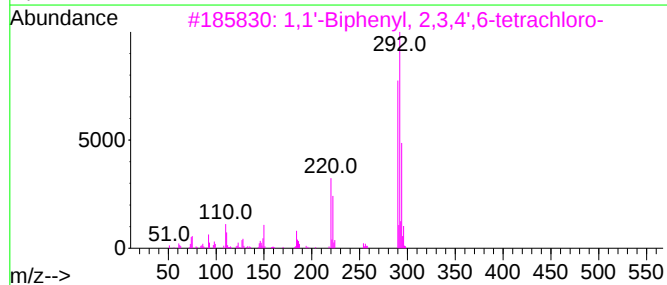
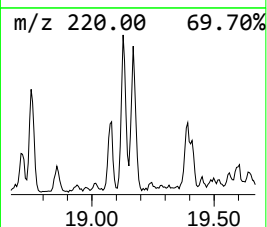
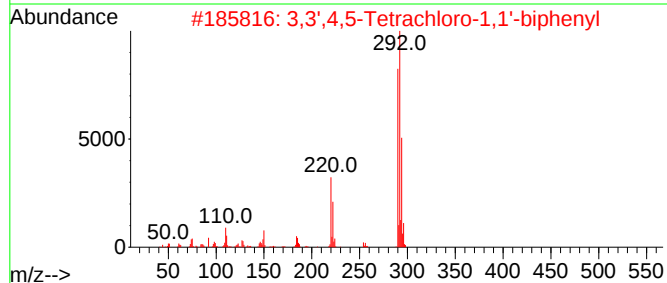
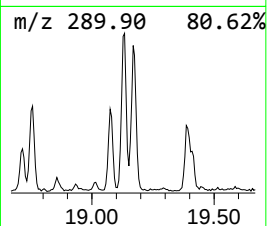
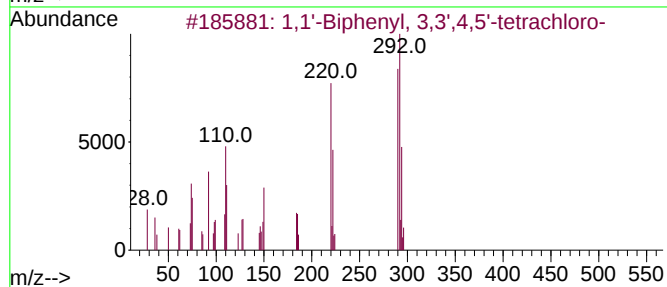
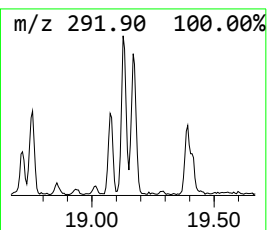
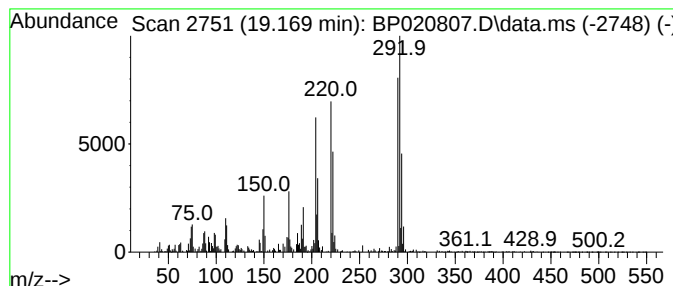
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1'-Biphenyl, 3,3',4,5'-te... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.169	3.23 ng/ul	431704	Phenanthrene-d10	17.169

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 3,3',4,5'-tetrach...	290	C12H6Cl4	041464-48-6	93		
2	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	93		
3	1,1'-Biphenyl, 2,3,4',6-tetrachl...	290	C12H6Cl4	052663-58-8	93		
4	1,1'-Biphenyl, 3,3',5,5'-tetrach...	290	C12H6Cl4	033284-52-5	93		
5	1,1'-Biphenyl, 2,3,5,6-tetrachloro-	290	C12H6Cl4	033284-54-7	93		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020807.D
Acq On : 06 Jul 2024 04:50
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Instrument :
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ClientSampleId :
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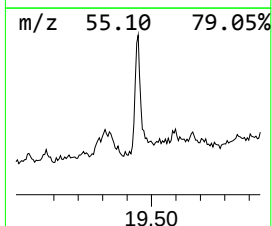
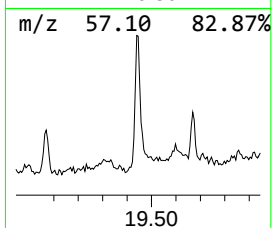
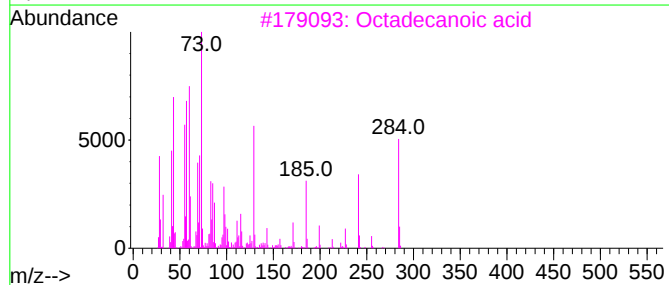
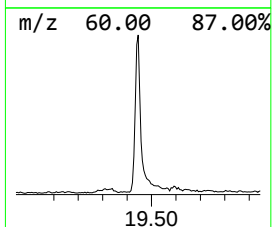
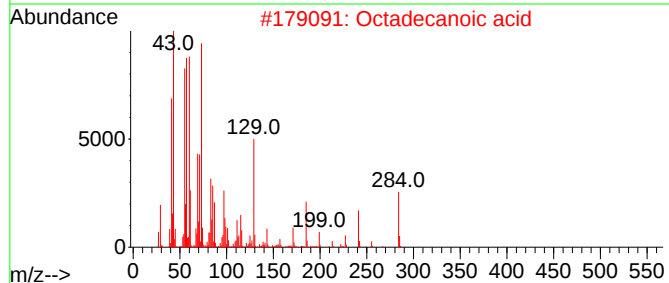
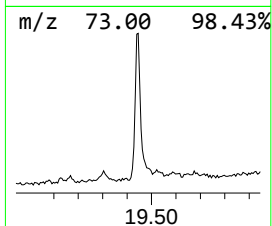
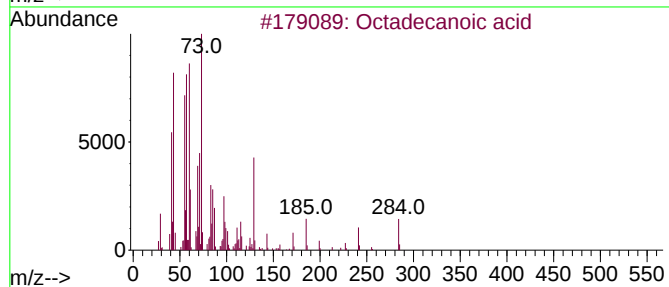
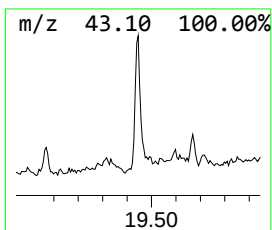
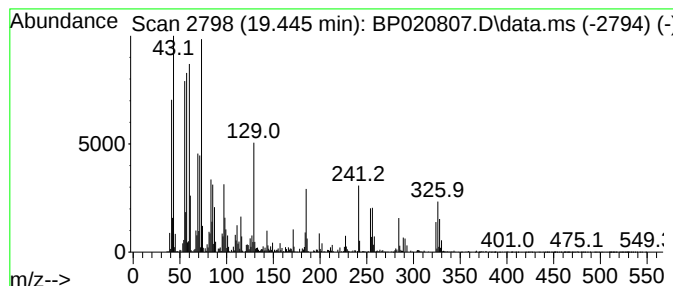
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	3.68 ng/ul	854948	Chrysene-d12	21.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	97
2			Octadecanoic acid	284	C18H36O2	000057-11-4	97
3			Octadecanoic acid	284	C18H36O2	000057-11-4	96
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	89
5			Heneicosanoic acid	326	C21H42O2	002363-71-5	78



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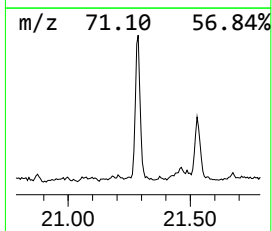
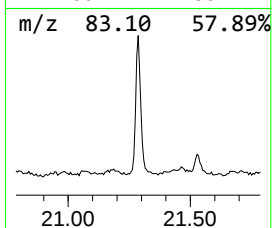
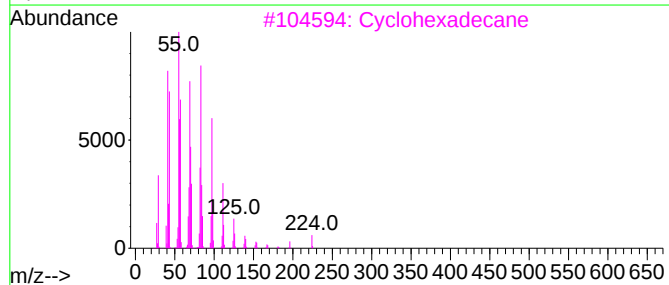
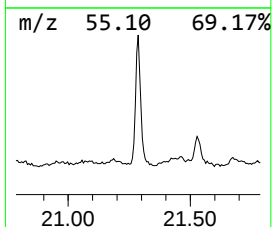
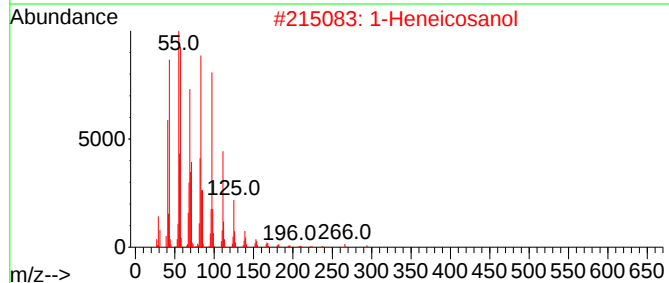
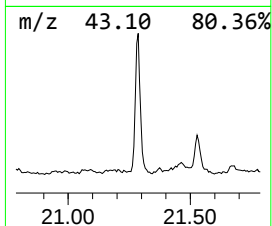
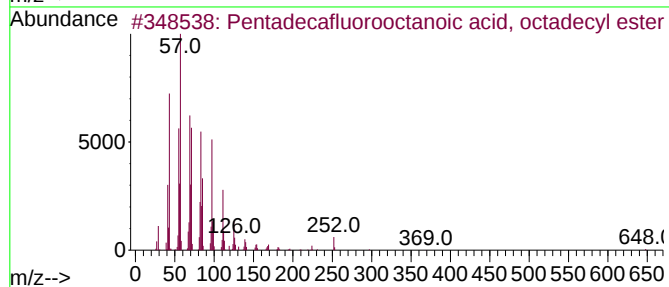
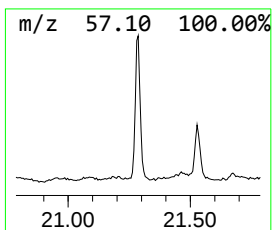
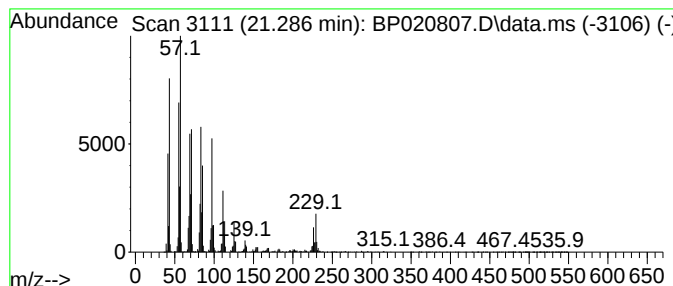
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Pentadecafluorooctanoic aci... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.286	8.50 ng/ul	1974170	Chrysene-d12	21.622	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2	1-Heneicosanol	312	C21H44O	015594-90-8	90
3	Cyclohexadecane	224	C16H32	000295-65-8	90
4	Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	89
5	Docosyl pentafluoropropionate	472	C25H45F5O2	1000351-80-9	87



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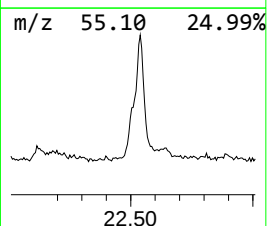
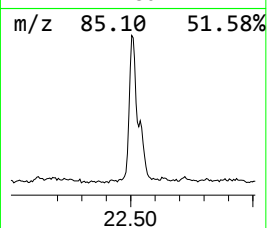
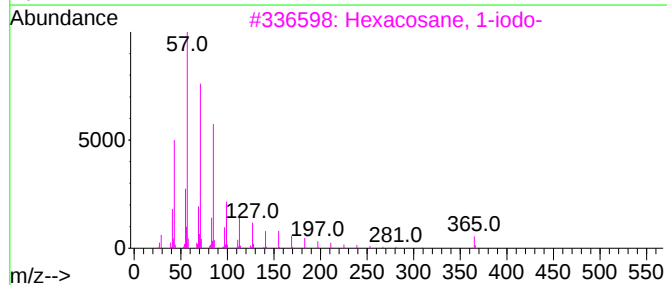
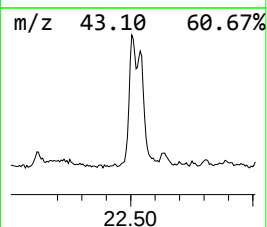
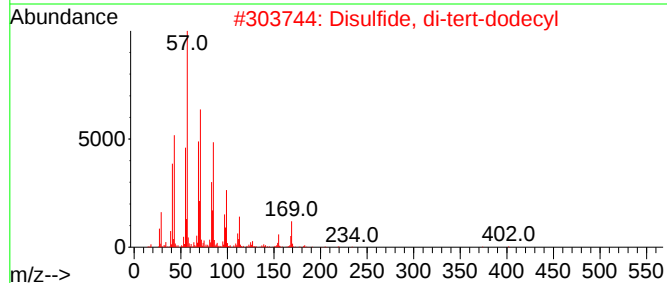
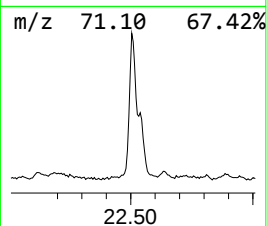
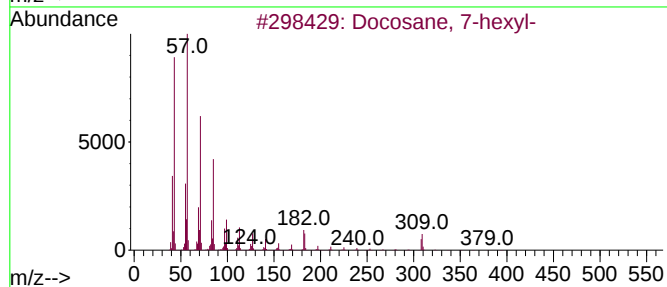
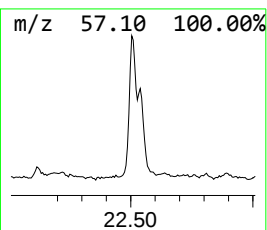
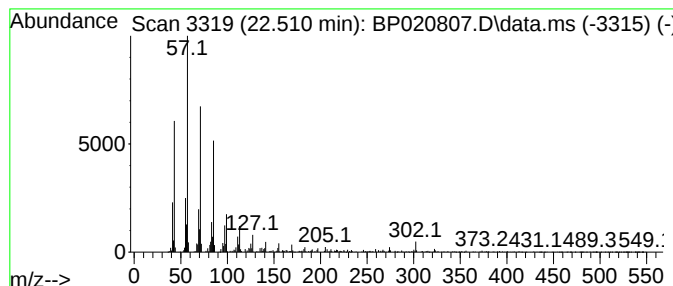
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.510	3.66 ng/ul	850506	Chrysene-d12	21.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
2			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	93
3			Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	90
4			Tetratetracontane	619	C44H90	007098-22-8	90
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020807.D
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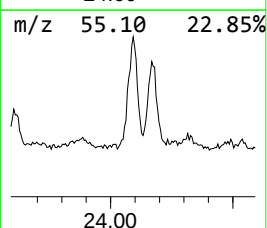
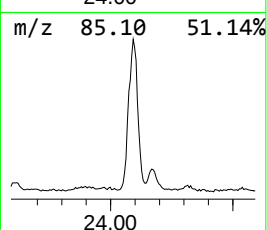
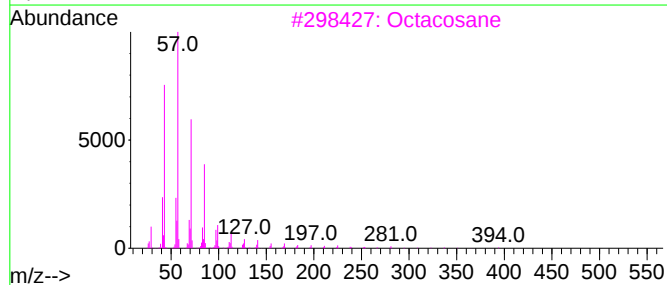
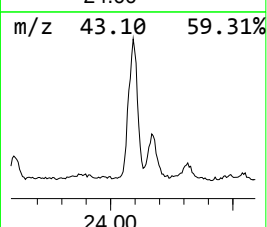
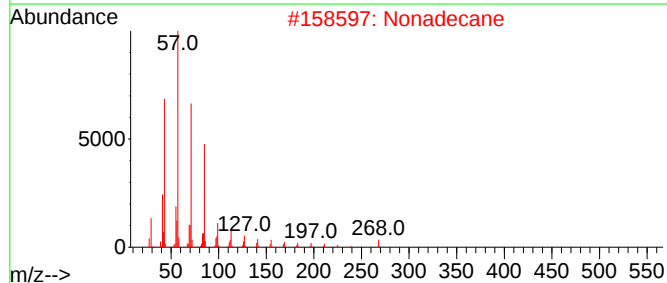
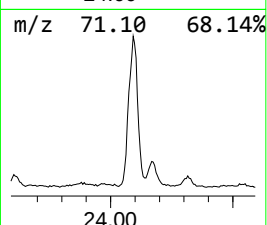
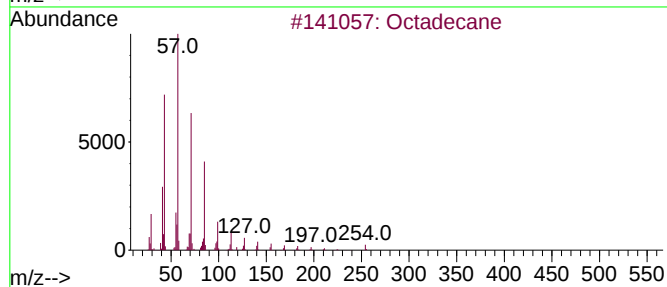
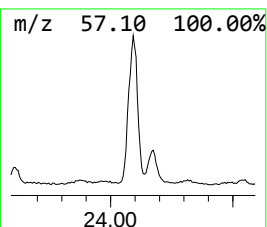
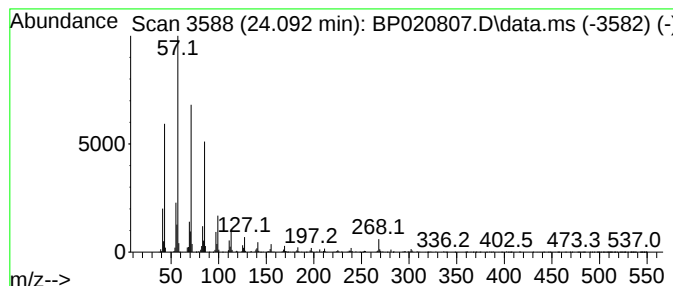
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.092	16.69 ng/ul	2238850	Perylene-d12	24.968

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Nonadecane	268	C19H40	000629-92-5	93
3			Octacosane	394	C28H58	000630-02-4	90
4			Tridecane, 6-propyl-	226	C16H34	055045-10-8	90
5			triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020807.D
Acq On : 06 Jul 2024 04:50
Operator : MA/JU
Sample : P2964-12
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CS6

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	4.881	4.2	ng/ul	239886	1	7.722	1155380	20.0
1,1'-Biphenyl, ...	17.792	3.5	ng/ul	461687	4	17.169	2676690	20.0
Anthracene, 2-m...	18.069	3.3	ng/ul	435846	4	17.169	2676690	20.0
n-Hexadecanoic ...	18.110	12.3	ng/ul	1641910	4	17.169	2676690	20.0
1,1'-Biphenyl, ...	18.275	4.8	ng/ul	648117	4	17.169	2676690	20.0
1,1'-Biphenyl, ...	19.169	3.2	ng/ul	431704	4	17.169	2676690	20.0
Octadecanoic acid	19.445	3.7	ng/ul	854948	5	21.622	4646260	20.0
Pentadecafluoro...	21.286	8.5	ng/ul	1974170	5	21.622	4646260	20.0
(DEL) Alkane: S...	22.510	3.7	ng/ul	850506	5	21.622	4646260	20.0
(DEL) Alkane: S...	24.092	16.7	ng/ul	2238850	6	24.968	2683630	20.0