

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
 Data File : BP020808.D
 Acq On : 06 Jul 2024 05:33
 Operator : MA/JU
 Sample : P2964-11
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4CS5

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: BP020808.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	44	47	58	rVB	35951	51660	1.29%	0.084%
2	3.540	90	94	108	rVB	83333	132731	3.32%	0.217%
3	3.775	130	134	139	rVB	84967	107085	2.68%	0.175%
4	3.899	151	155	161	rVV8	8069	15884	0.40%	0.026%
5	3.987	165	170	180	rVB	110900	159947	4.00%	0.261%
6	4.146	193	197	203	rBV5	10339	19639	0.49%	0.032%
7	4.352	229	232	239	rVB5	11931	16480	0.41%	0.027%
8	4.487	251	255	259	rBV	13332	19294	0.48%	0.032%
9	4.534	259	263	268	rVV	40235	60000	1.50%	0.098%
10	4.587	268	272	279	rVB	36212	50464	1.26%	0.082%
11	4.887	317	323	331	rBV	182154	261029	6.54%	0.426%
12	4.987	335	340	347	rVB2	17691	31427	0.79%	0.051%
13	5.269	382	388	395	rVB6	10021	23688	0.59%	0.039%
14	5.405	407	411	419	rVB	12738	24214	0.61%	0.040%
15	5.763	466	472	477	rBV3	28549	46173	1.16%	0.075%
16	5.846	482	486	495	rVB	21066	39537	0.99%	0.065%
17	6.381	567	577	578	rBV9	9094	17852	0.45%	0.029%
18	6.769	638	643	648	rVV3	18085	33350	0.83%	0.054%
19	6.922	658	669	681	rVV	602220	1090922	27.31%	1.782%
20	7.075	688	695	701	rBV	543753	809720	20.27%	1.322%
21	7.269	722	728	736	rBV	665570	1118190	27.99%	1.826%
22	7.340	736	740	752	rVB4	61072	133911	3.35%	0.219%
23	7.728	792	806	814	rBV	880311	1394924	34.92%	2.278%
24	7.999	848	852	858	rVB9	8355	16892	0.42%	0.028%
25	8.257	884	896	900	rBV4	15786	33865	0.85%	0.055%
26	8.310	900	905	909	rVV7	7006	15911	0.40%	0.026%
27	8.357	909	913	919	rVB6	11833	22527	0.56%	0.037%
28	8.434	919	926	941	rBV	787985	1336346	33.46%	2.182%
29	8.546	941	945	952	rVV	37463	63719	1.60%	0.104%
30	8.804	982	989	994	rBV8	14430	29883	0.75%	0.049%
31	8.875	994	1001	1007	rBV	637774	1089601	27.28%	1.780%
32	9.028	1018	1027	1036	rVB	12166	44668	1.12%	0.073%
33	9.434	1083	1096	1102	rBV	61824	116952	2.93%	0.191%
34	9.587	1115	1122	1133	rBV	556757	944514	23.65%	1.543%
35	9.834	1160	1164	1171	rVB5	17666	34185	0.86%	0.056%
36	10.134	1207	1215	1233	rBV	741256	1425164	35.68%	2.328%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
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37	10.281	1235	1240	1246	rVB9	7502	18622	0.47%	0.030%
38	10.498	1264	1277	1283	rBV	1182894	2018991	50.55%	3.297%
39	10.551	1283	1286	1292	rVV	111153	172200	4.31%	0.281%
40	10.645	1296	1302	1321	rVB	115481	270908	6.78%	0.442%
41	11.034	1361	1368	1375	rBV4	27608	49102	1.23%	0.080%
42	11.245	1396	1404	1411	rBV2	133776	293950	7.36%	0.480%
43	11.298	1411	1413	1418	rVB5	9617	16815	0.42%	0.027%
44	11.498	1443	1447	1452	rBV	12400	20241	0.51%	0.033%
45	11.910	1511	1517	1522	rVB	27786	46058	1.15%	0.075%
46	12.010	1528	1534	1539	rBV4	14963	32427	0.81%	0.053%
47	12.081	1542	1546	1552	rVV	18212	33324	0.83%	0.054%
48	12.151	1553	1558	1567	rVB	64171	116994	2.93%	0.191%
49	12.298	1576	1583	1588	rBV6	10953	27812	0.70%	0.045%
50	12.387	1592	1598	1605	rVB	52608	77497	1.94%	0.127%
51	12.634	1629	1640	1646	rVV	9010	24124	0.60%	0.039%
52	12.869	1676	1680	1686	rVB5	13573	22494	0.56%	0.037%
53	13.163	1724	1730	1737	rBV2	47825	93160	2.33%	0.152%
54	13.381	1762	1767	1770	rBV2	14075	25381	0.64%	0.041%
55	13.528	1783	1792	1797	rBV4	32809	93286	2.34%	0.152%
56	13.669	1810	1816	1822	rBV2	48478	92663	2.32%	0.151%
57	13.728	1822	1826	1828	rBV2	32345	51992	1.30%	0.085%
58	13.769	1828	1833	1839	rVV	1462125	1928753	48.29%	3.150%
59	13.904	1853	1856	1859	rVB2	24206	29508	0.74%	0.048%
60	14.045	1869	1880	1895	rBV2	1451796	2849241	71.33%	4.653%
61	14.257	1908	1916	1920	rBV2	38731	57187	1.43%	0.093%
62	14.357	1924	1933	1939	rBV	1698321	2637198	66.02%	4.307%
63	14.416	1939	1943	1951	rVB2	78518	148571	3.72%	0.243%
64	14.492	1951	1956	1961	rBV4	9954	19683	0.49%	0.032%
65	14.598	1963	1974	1983	rBV	477435	955496	23.92%	1.561%
66	14.757	1993	2001	2010	rVB3	109611	289178	7.24%	0.472%
67	14.834	2010	2014	2020	rVB	26677	41910	1.05%	0.068%
68	14.986	2036	2040	2043	rVV2	26587	34159	0.86%	0.056%
69	15.033	2045	2048	2052	rVB	30384	38590	0.97%	0.063%
70	15.157	2065	2069	2073	rVV6	33162	52723	1.32%	0.086%
71	15.222	2077	2080	2082	rVV	22012	23663	0.59%	0.039%
72	15.369	2098	2105	2110	rBV	2078658	3053843	76.46%	4.987%
73	15.416	2110	2113	2119	rVV2	88389	140659	3.52%	0.230%
74	15.504	2123	2128	2133	rBV	459079	624096	15.62%	1.019%
75	15.633	2147	2150	2157	rVB6	28307	50017	1.25%	0.082%
76	15.728	2162	2166	2172	rVB	56069	95412	2.39%	0.156%

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77	15.875	2187	2191	2198	rVB3	47637	93172	2.33%	0.152%
78	16.122	2226	2233	2237	rBV3	129042	262324	6.57%	0.428%
79	16.557	2304	2307	2312	rBV5	50608	64179	1.61%	0.105%
80	16.816	2347	2351	2356	rBV2	112438	170105	4.26%	0.278%
81	16.916	2366	2368	2372	rVB	73146	69994	1.75%	0.114%
82	16.980	2375	2379	2386	rVB2	119635	203853	5.10%	0.333%
83	17.175	2405	2412	2415	rBV	1843084	2994648	74.97%	4.891%
84	17.216	2415	2419	2423	rVV	1547456	2115765	52.97%	3.455%
85	17.269	2423	2428	2432	rVV	2131140	3147375	78.80%	5.140%
86	17.304	2432	2434	2439	rVB	389146	419072	10.49%	0.684%
87	17.357	2440	2443	2449	rBV4	69576	137512	3.44%	0.225%
88	17.786	2511	2516	2522	rVB3	168022	319817	8.01%	0.522%
89	17.910	2534	2537	2543	rVB3	125489	205091	5.13%	0.335%
90	18.110	2562	2571	2579	rBV2	1025960	2577132	64.52%	4.209%
91	18.274	2594	2599	2603	rBV2	952139	1537114	38.48%	2.510%
92	18.380	2615	2617	2625	rVB4	268825	442460	11.08%	0.723%
93	18.574	2647	2650	2652	rBV	340169	408516	10.23%	0.667%
94	19.133	2742	2745	2748	rBV2	542533	601451	15.06%	0.982%
95	19.310	2770	2775	2780	rBV	2660143	3994254	100.00%	6.523%
96	19.445	2794	2798	2804	rBV3	801568	1612563	40.37%	2.634%
97	19.663	2830	2835	2837	rBV2	2444902	3527305	88.31%	5.761%
98	19.692	2837	2840	2848	rVB	1952566	3008378	75.32%	4.913%
99	21.298	3110	3113	3120	rVB3	1422819	2314859	57.95%	3.781%
100	21.627	3165	3169	3175	rVB3	1768245	3502847	87.70%	5.721%

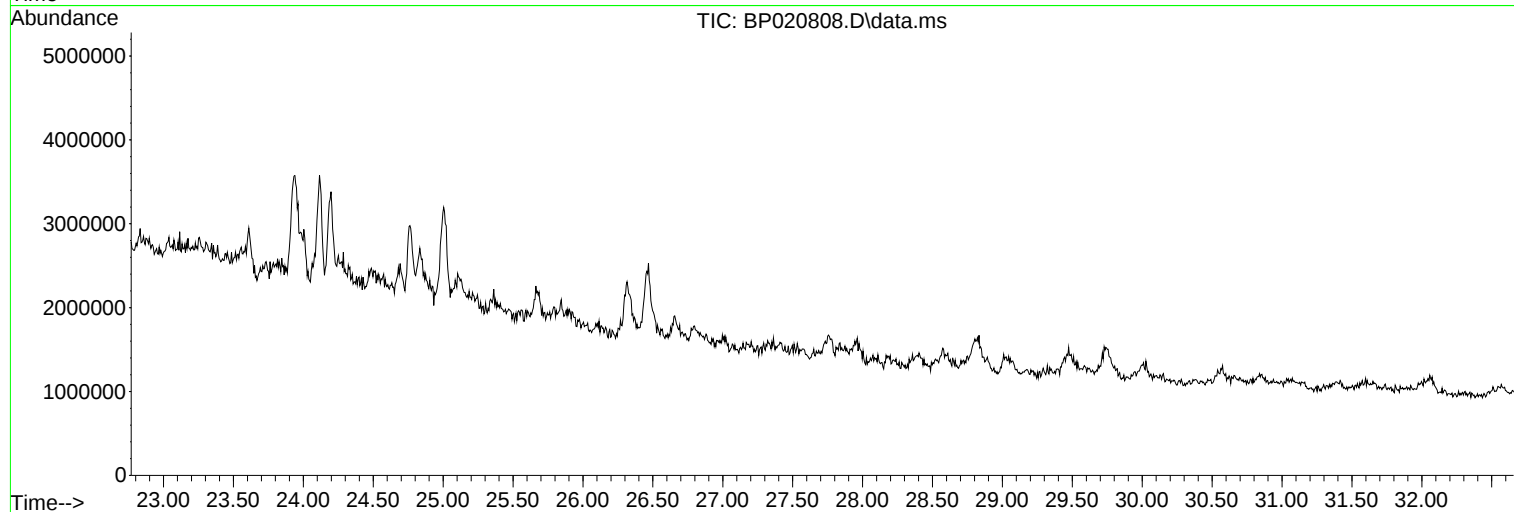
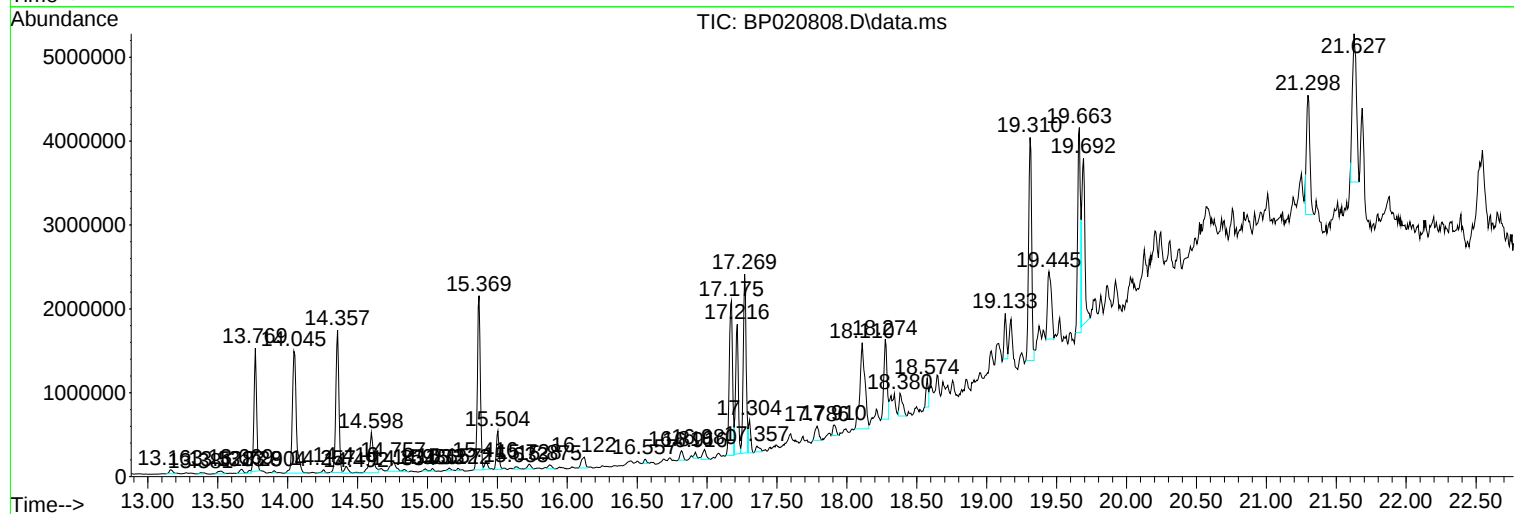
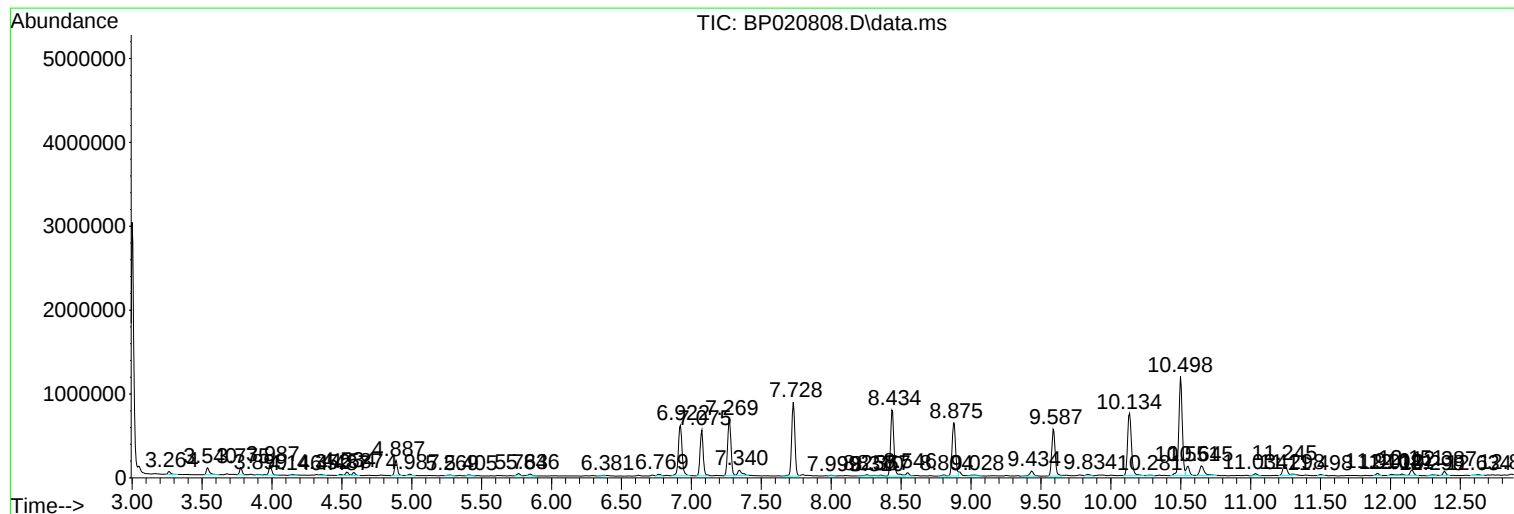
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Instrument :
BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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



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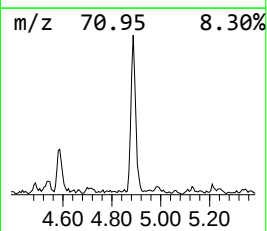
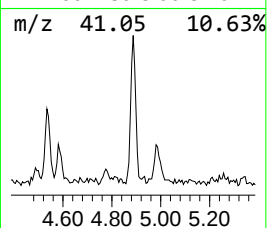
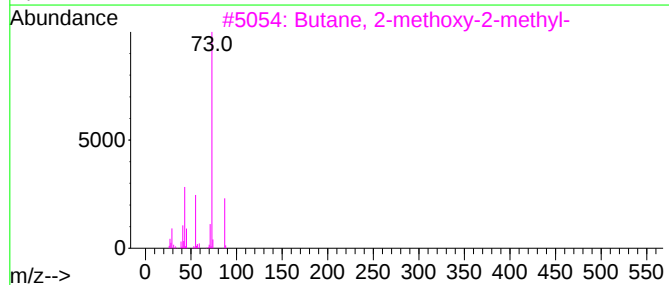
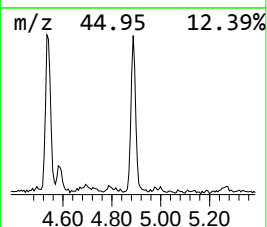
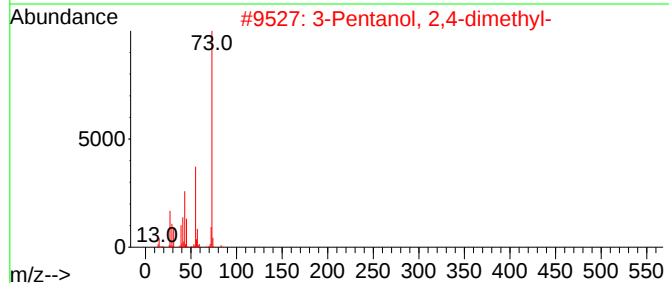
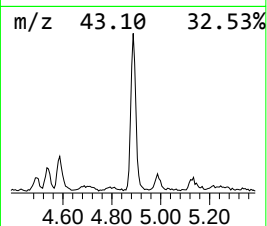
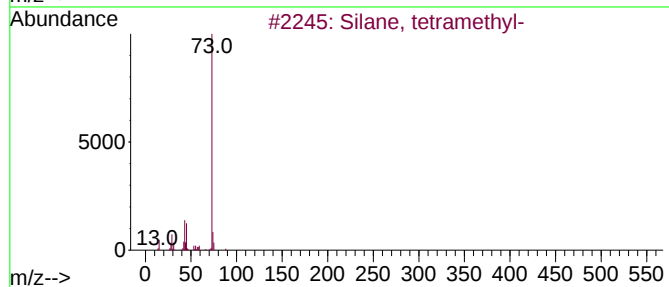
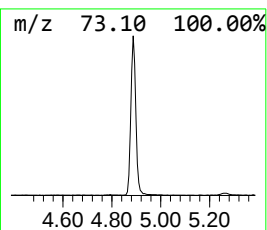
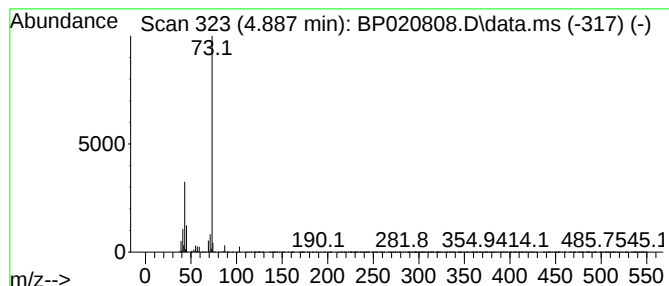
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.887	3.74 ng/ul	261029	1,4-Dichlorobenzene-d4	7.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, tetramethyl-	88	C4H12Si	000075-76-3	47
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	36
3			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	36
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	33
5			1,3-Dioxolane, 2-(6-heptynyl)-	168	C10H16O2	025595-82-8	28



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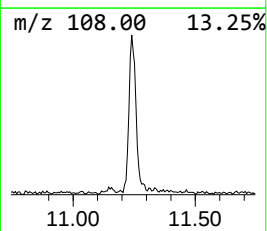
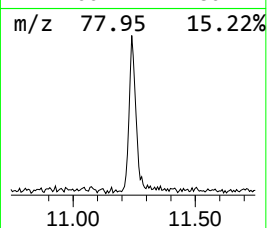
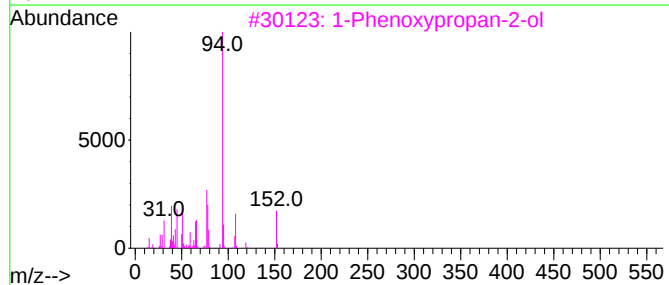
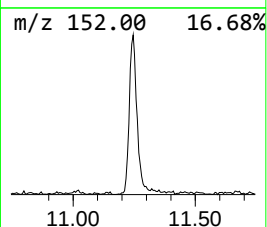
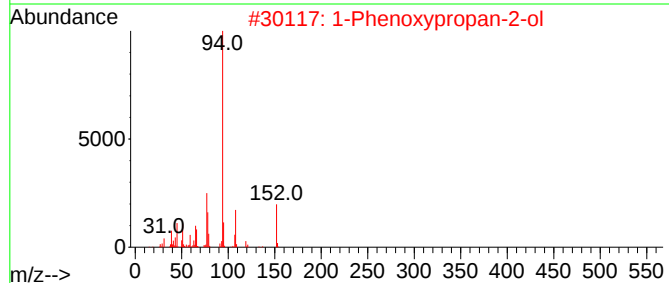
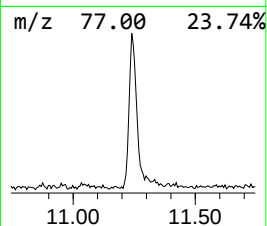
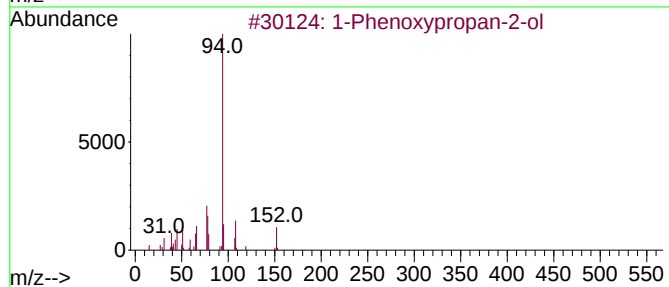
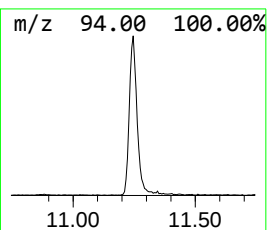
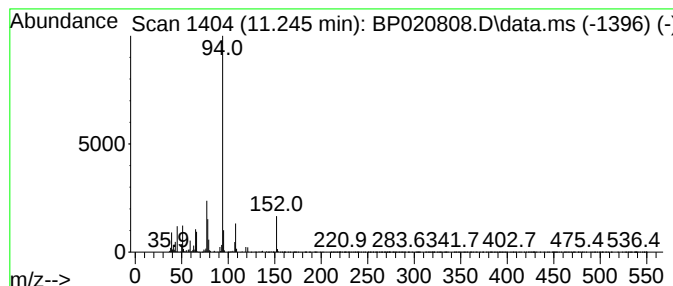
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062524.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.246	2.91 ng/ul	293950	Naphthalene-d8	10.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83



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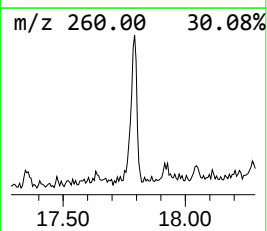
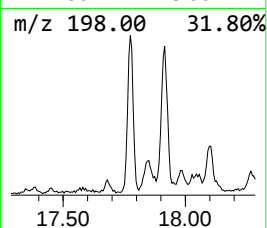
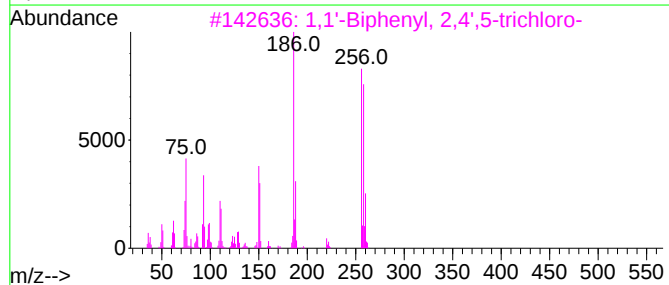
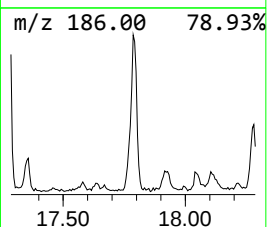
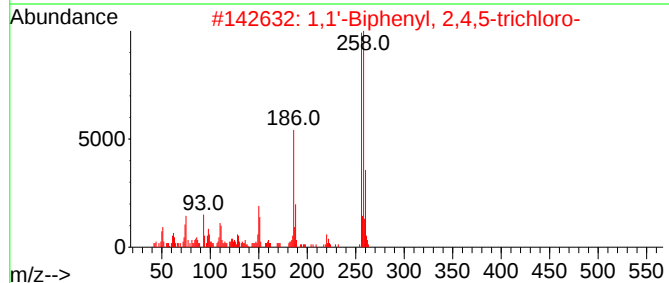
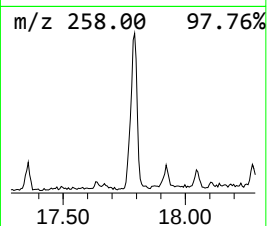
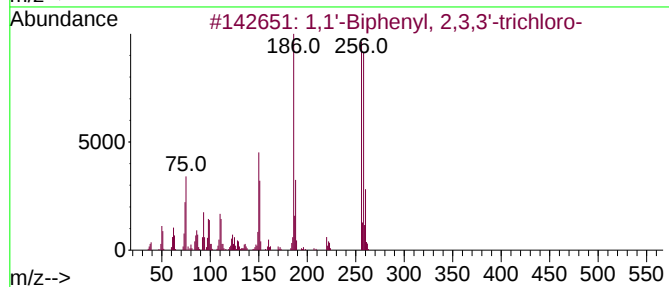
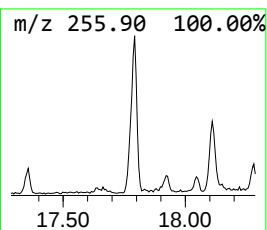
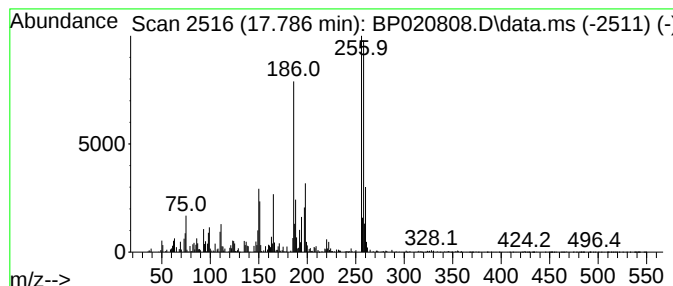
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.786	2.14 ng/ul	319817	Phenanthrene-d10	17.175

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, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	97		
3	1,1'-Biphenyl, 2,4',5-trichloro-	256	C12H7Cl3	016606-02-3	96		
4	1,1'-Biphenyl, 2,4,5-trichloro-	256	C12H7Cl3	015862-07-4	95		
5	1,1'-Biphenyl, 2,3,4'-Trichloro-	256	C12H7Cl3	038444-85-8	95		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
Operator : MA/JU
Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CS5

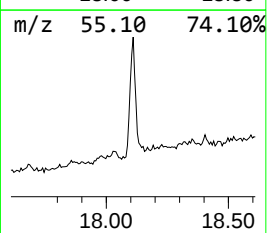
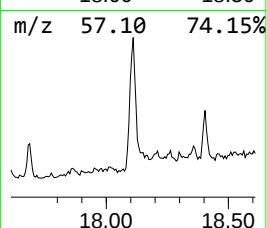
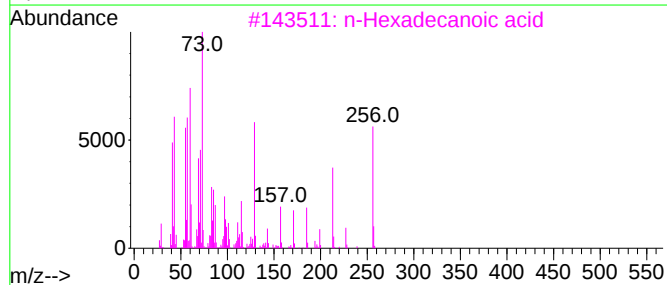
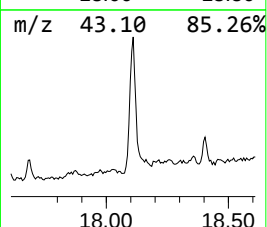
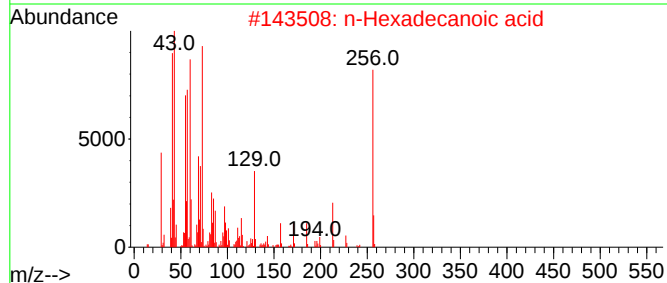
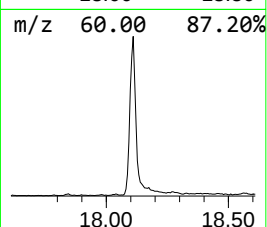
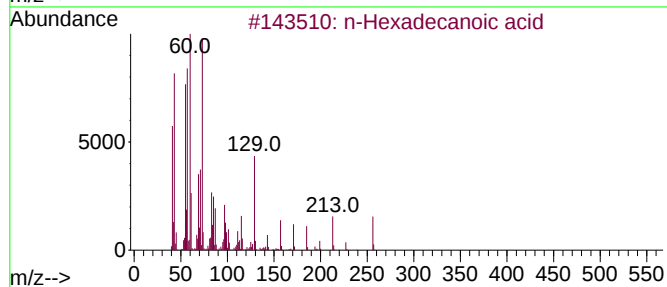
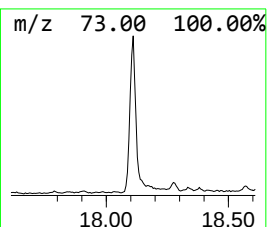
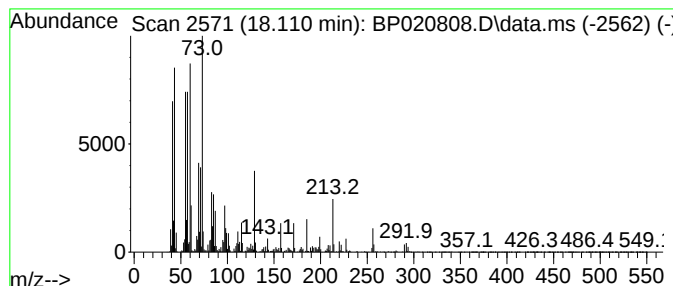
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.110	17.21 ng/ul	2577130	Phenanthrene-d10	17.175

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	98
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
Operator : MA/JU
Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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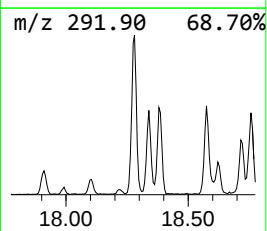
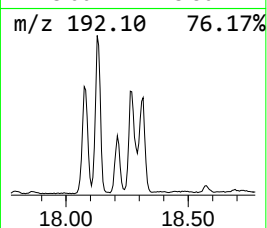
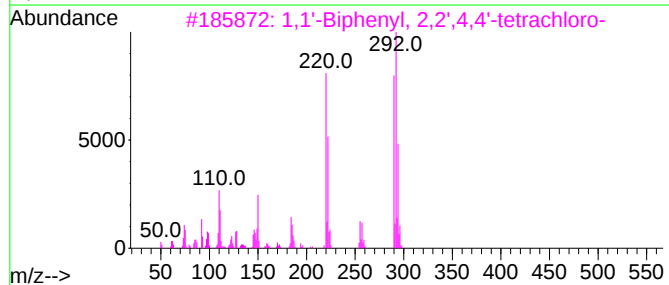
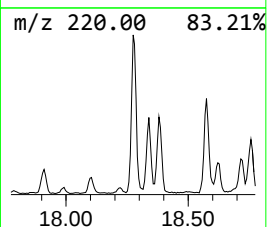
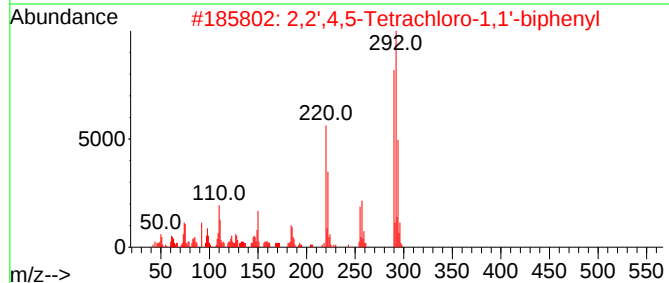
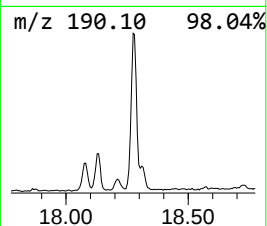
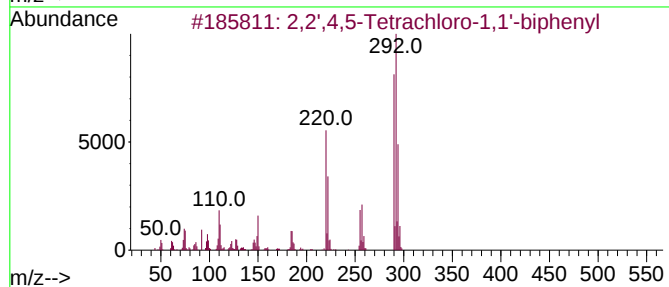
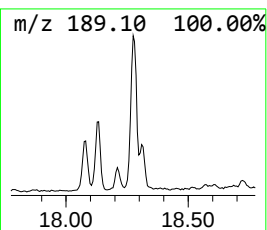
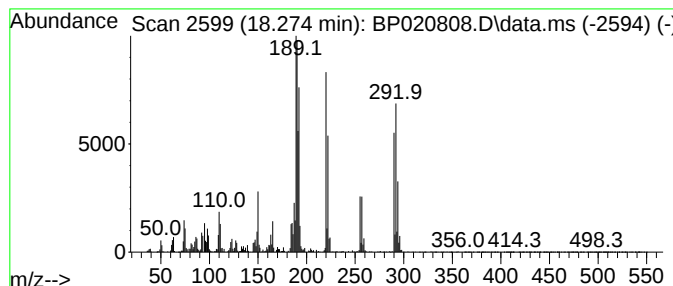
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2,2',4,5-Tetrachloro-1,1'-b... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.275	10.27 ng/ul	1537110	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	96		
2	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	96		
3	1,1'-Biphenyl, 2,2',4,4'-tetrachloro-	290	C12H6Cl4	002437-79-8	96		
4	1,1'-Biphenyl, 2,2',4,5'-tetrachloro-	290	C12H6Cl4	041464-40-8	96		
5	1,1'-Biphenyl, 2,2',6,6'-tetrachloro-	290	C12H6Cl4	015968-05-5	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
Operator : MA/JU
Sample : P2964-11
Misc :
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ClientSampleId :
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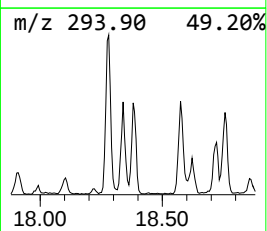
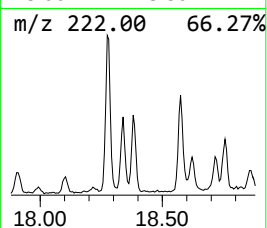
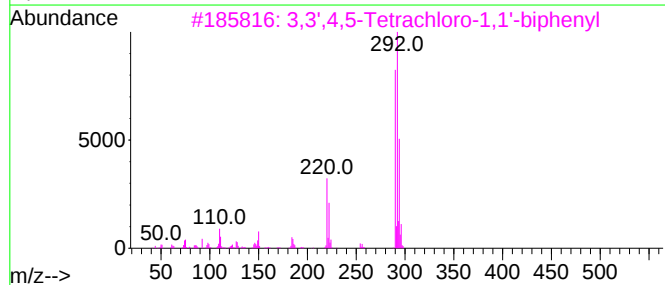
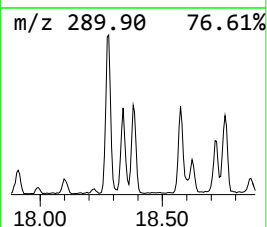
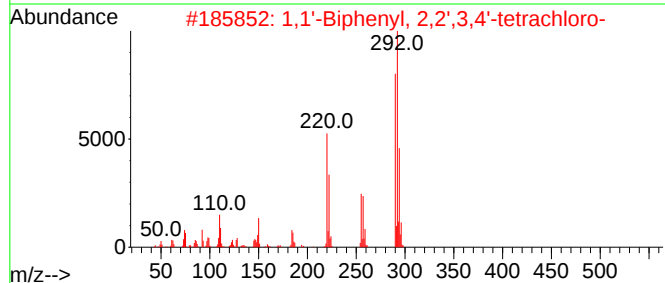
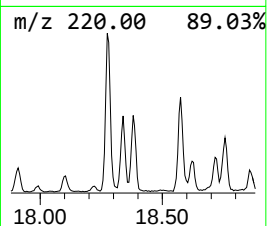
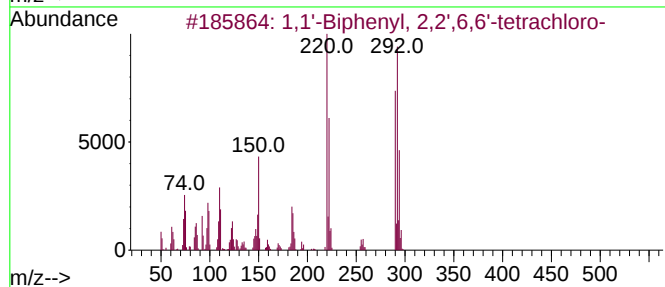
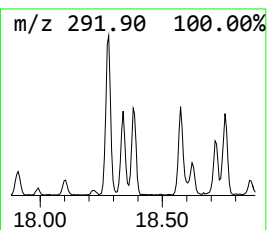
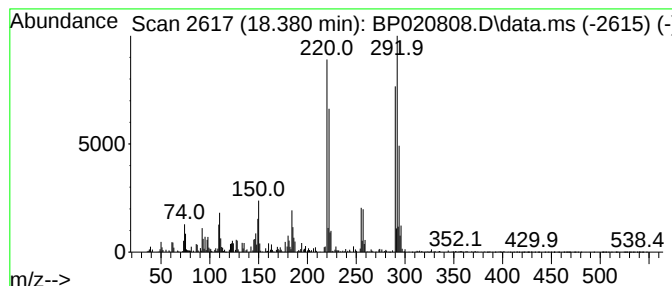
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,2',6,6'-te... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.380	2.96 ng/ul	442460	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99		
2	1,1'-Biphenyl, 2,2',3,4'-tetrach...	290	C12H6Cl4	036559-22-5	96		
3	3,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-49-1	96		
4	2,3',5',6-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	074338-23-1	96		
5	2,3',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	073575-53-8	96		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
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Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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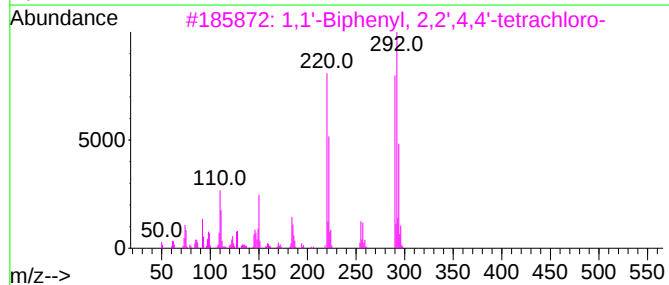
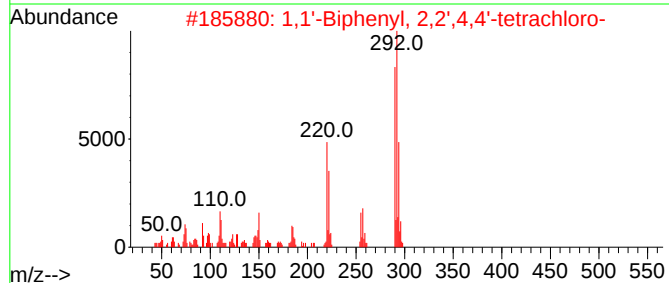
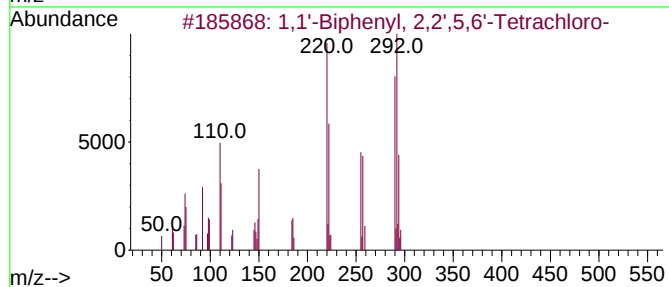
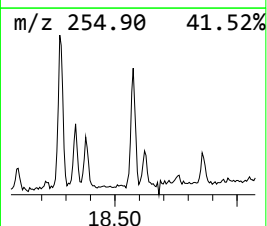
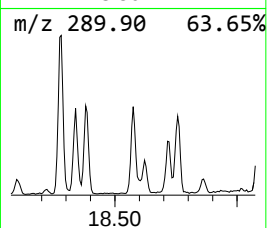
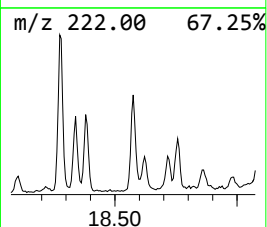
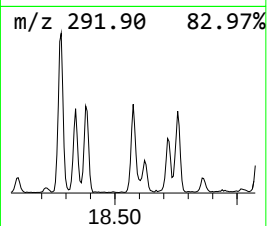
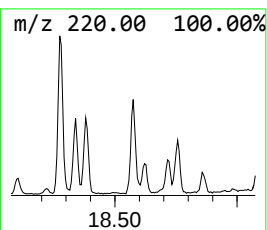
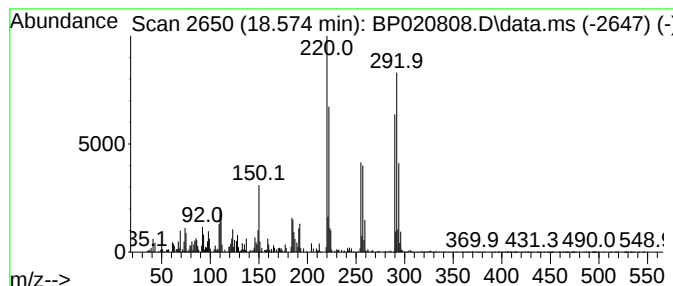
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,2',5,6'-Te... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.575	2.73 ng/ul	408516	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',5,6'-Tetrach...	290	C12H6Cl4	041464-41-9	99		
2	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
3	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
4	1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99		
5	2,2',4,5-Tetrachloro-1,1'-biphenyl	290	C12H6Cl4	070362-47-9	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
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Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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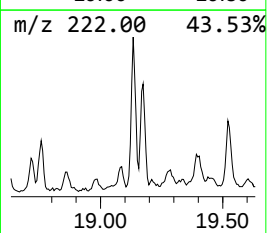
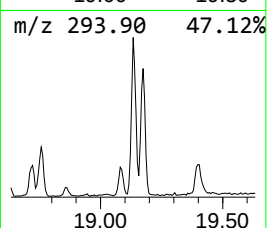
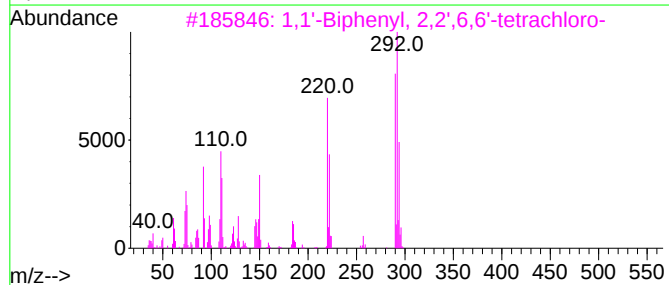
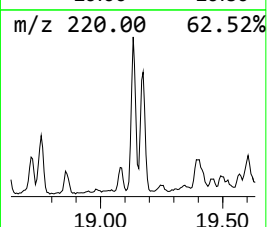
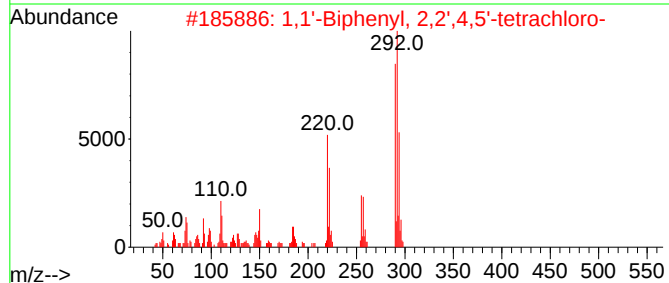
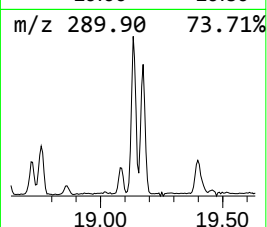
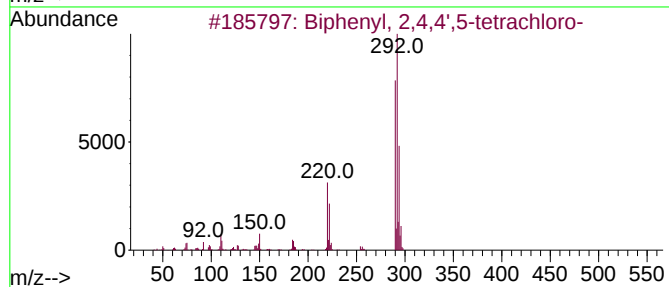
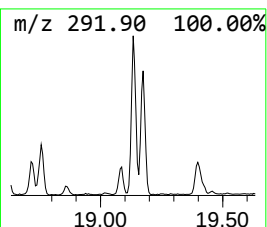
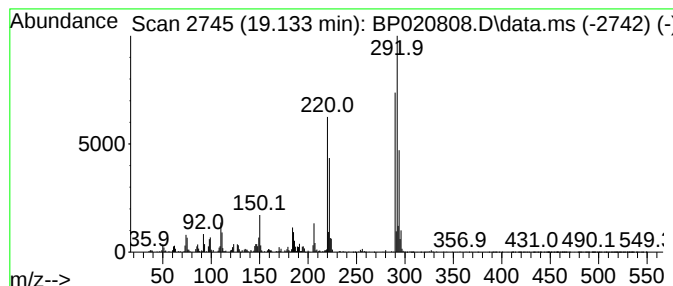
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Biphenyl, 2,4,4',5-tetrachl... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.133	4.02 ng/ul	601451	Phenanthrene-d10	17.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Biphenyl, 2,4,4',5-tetrachloro-	290	C12H6Cl4	032690-93-0	99
2			1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	99
3			1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99
4			1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	98
5			1,1'-Biphenyl, 2,2',4,5'-tetrach...	290	C12H6Cl4	041464-40-8	98



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
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Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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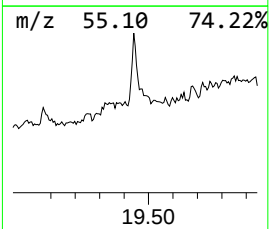
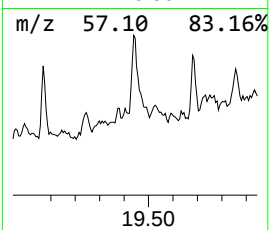
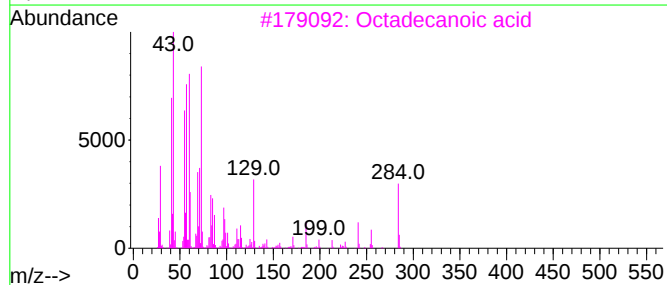
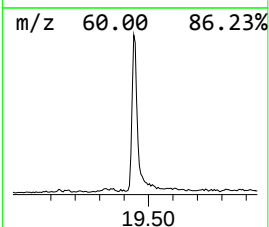
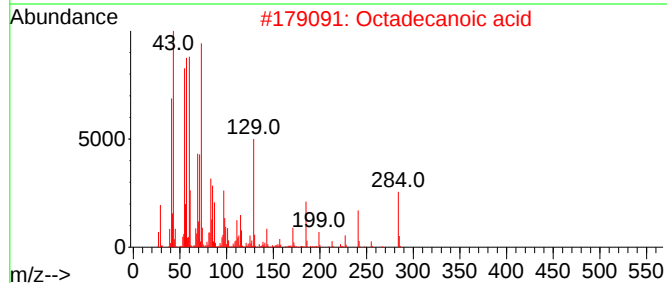
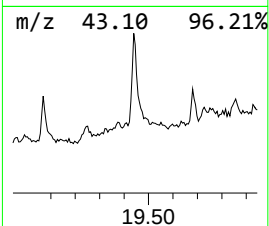
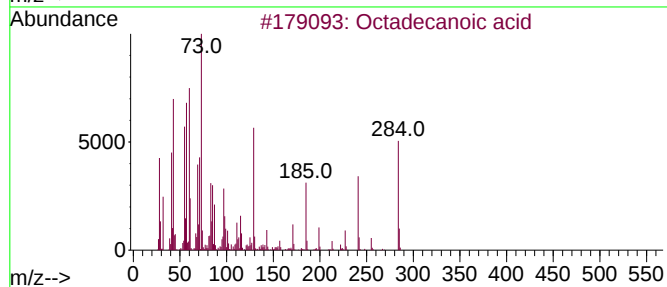
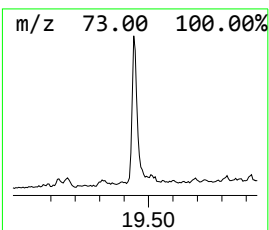
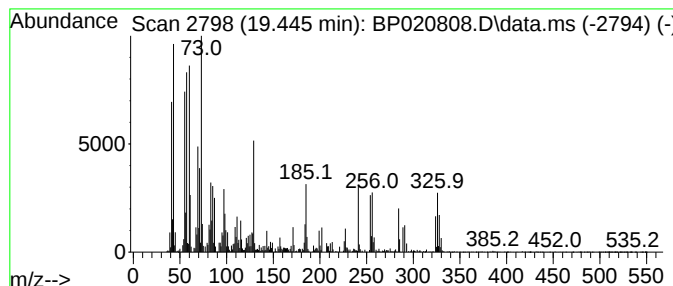
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	9.21 ng/ul	1612560	Chrysene-d12	21.633

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	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Octadecanoic acid	284	C18H36O2	000057-11-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
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Operator : MA/JU
Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

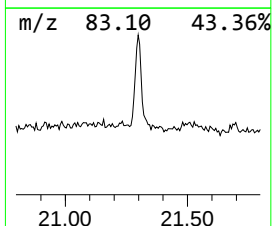
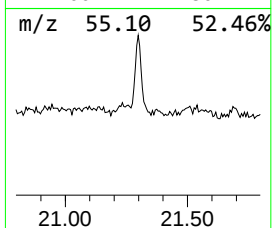
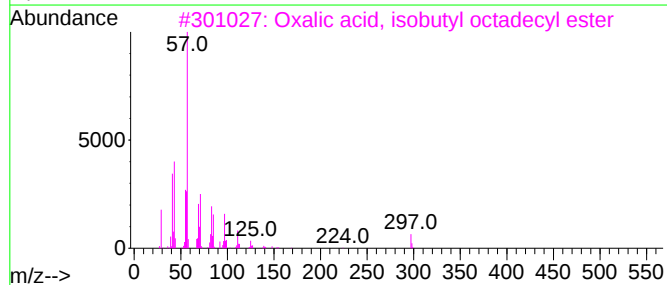
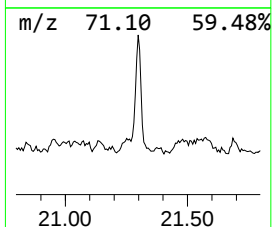
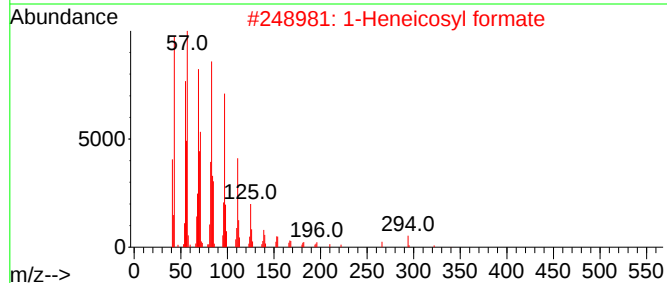
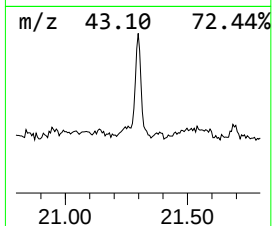
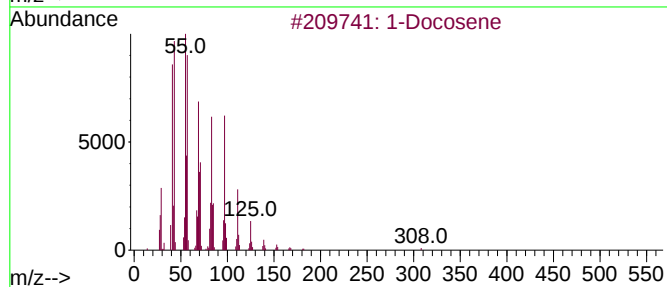
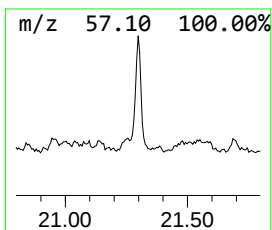
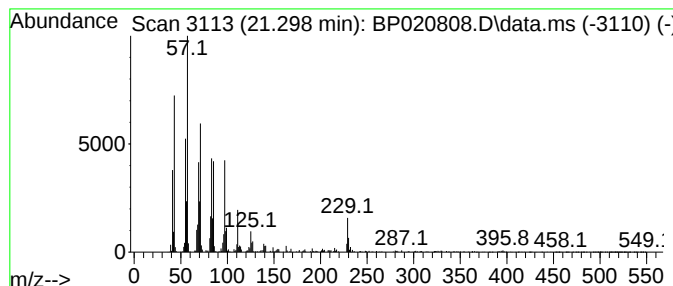
Instrument :
BNA_P
ClientSampleId :
A4CS5

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 11 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.298	13.22 ng/ul	2314860	Chrysene-d12	21.633	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	92
2	1-Heneicosyl formate	340	C22H44O2	077899-03-7	91
3	Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	87
4	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	87
5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070524\
Data File : BP020808.D
Acq On : 06 Jul 2024 05:33
Operator : MA/JU
Sample : P2964-11
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4CS5

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	4.887	3.7	ng/ul	261029	1	7.728	1394920	20.0
1-Phenoxypropan...	11.246	2.9	ng/ul	293950	2	10.498	2018990	20.0
1,1'-Biphenyl, ...	17.786	2.1	ng/ul	319817	4	17.175	2994650	20.0
n-Hexadecanoic ...	18.110	17.2	ng/ul	2577130	4	17.175	2994650	20.0
2,2',4,5-Tetrac...	18.275	10.3	ng/ul	1537110	4	17.175	2994650	20.0
1,1'-Biphenyl, ...	18.380	3.0	ng/ul	442460	4	17.175	2994650	20.0
1,1'-Biphenyl, ...	18.575	2.7	ng/ul	408516	4	17.175	2994650	20.0
Biphenyl, 2,4,4...	19.133	4.0	ng/ul	601451	4	17.175	2994650	20.0
Octadecanoic acid	19.445	9.2	ng/ul	1612560	5	21.633	3502850	20.0
(DEL) Alkane: S...	21.298	13.2	ng/ul	2314860	5	21.633	3502850	20.0