

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
 Data File : BP017074.D
 Acq On : 10 Sep 2023 04:23
 Operator : MA/JU
 Sample : 04227-11
 Misc :
 ALS Vial : 86 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBKT7

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-BP090523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017074.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.275	27	32	37	rBV	1342063	1881728	12.29%	0.909%
2	3.322	37	40	44	rVV	284244	401377	2.62%	0.194%
3	3.364	44	47	54	rVB	108136	151042	0.99%	0.073%
4	3.599	84	87	92	rBV	77682	98916	0.65%	0.048%
5	3.764	111	115	131	rBV	1252843	1870306	12.21%	0.903%
6	3.958	143	148	157	rVB	106047	166522	1.09%	0.080%
7	4.070	162	167	176	rVB	448664	628153	4.10%	0.303%
8	4.599	250	257	276	rVB	6387446	9411729	61.45%	4.546%
9	5.234	358	365	376	rBV	4566960	6724071	43.90%	3.248%
10	5.417	390	396	402	rBV	54926	90198	0.59%	0.044%
11	5.605	424	428	433	rVB2	72203	119320	0.78%	0.058%
12	5.793	455	460	466	rBV2	50358	98770	0.64%	0.048%
13	5.928	473	483	489	rBV	188046	307417	2.01%	0.148%
14	6.216	528	532	545	rVB6	40064	92261	0.60%	0.045%
15	6.405	554	564	571	rBV	171570	316286	2.07%	0.153%
16	6.487	573	578	582	rVV	87179	124864	0.82%	0.060%
17	6.658	602	607	613	rVV	97255	131440	0.86%	0.063%
18	6.905	643	649	655	rBV2	47432	91337	0.60%	0.044%
19	7.122	679	686	701	rVB	885749	1439779	9.40%	0.695%
20	7.311	709	718	729	rBV	3297731	5081443	33.18%	2.455%
21	7.487	741	748	754	rBV	3073885	4950322	32.32%	2.391%
22	7.528	754	755	761	rVV	232788	249866	1.63%	0.121%
23	7.593	761	766	770	rVV4	42375	89470	0.58%	0.043%
24	7.646	770	775	777	rVV5	52751	102343	0.67%	0.049%
25	7.675	778	780	789	rVB2	71103	108140	0.71%	0.052%
26	7.799	794	801	803	rVV	63704	99151	0.65%	0.048%
27	7.840	803	808	814	rVV2	278569	562534	3.67%	0.272%
28	7.899	814	818	823	rVV	107235	161249	1.05%	0.078%
29	7.969	823	830	834	rVV	2294716	3781873	24.69%	1.827%
30	8.005	834	836	842	rVV	801187	1014423	6.62%	0.490%
31	8.246	870	877	881	rBV4	65942	112820	0.74%	0.054%
32	8.316	882	889	899	rVB	5978959	9293321	60.68%	4.489%
33	8.428	899	908	911	rVB6	42342	101657	0.66%	0.049%
34	8.587	927	935	941	rBV2	69123	120799	0.79%	0.058%
35	8.681	943	951	961	rBV	2635998	4213945	27.51%	2.036%
36	8.769	961	966	971	rVV3	54503	125578	0.82%	0.061%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
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37	8.946	992	996	1003	rBV	49988	86664	0.57%	0.042%
38	9.169	1025	1034	1044	rVB	3752435	6172830	40.30%	2.982%
39	9.287	1044	1054	1058	rBV6	51151	129028	0.84%	0.062%
40	9.334	1058	1062	1066	rVV3	75802	112184	0.73%	0.054%

41	9.381	1066	1070	1077	rVB2	142287	248964	1.63%	0.120%
42	9.752	1128	1133	1138	rVV	128890	199181	1.30%	0.096%
43	9.822	1138	1145	1149	rVV4	83683	169962	1.11%	0.082%
44	9.881	1149	1155	1162	rVV	3055616	4896341	31.97%	2.365%
45	9.946	1162	1166	1170	rVV3	73233	126922	0.83%	0.061%

46	10.034	1170	1181	1189	rVV5	74627	205796	1.34%	0.099%
47	10.140	1195	1199	1206	rVV3	82011	226973	1.48%	0.110%
48	10.204	1206	1210	1216	rVV2	186159	312062	2.04%	0.151%
49	10.404	1235	1244	1266	rVB	4282296	7393459	48.27%	3.571%
50	10.646	1276	1285	1290	rBV	834678	1432526	9.35%	0.692%

51	10.693	1290	1293	1299	rVV2	43413	89534	0.58%	0.043%
52	10.793	1303	1310	1318	rVV	3169182	5271491	34.42%	2.546%
53	10.875	1321	1324	1331	rVV3	55561	100497	0.66%	0.049%
54	10.957	1331	1338	1351	rVV	3285334	5512369	35.99%	2.663%
55	11.063	1352	1356	1364	rVB6	47133	109963	0.72%	0.053%

56	11.169	1369	1374	1380	rVB	100819	180695	1.18%	0.087%
57	11.546	1431	1438	1441	rBV5	40211	85542	0.56%	0.041%
58	11.875	1488	1494	1497	rBV2	145463	218208	1.42%	0.105%
59	12.063	1520	1526	1533	rBV3	76291	164508	1.07%	0.079%
60	12.222	1543	1553	1559	rVV5	194871	534616	3.49%	0.258%

61	12.375	1572	1579	1585	rVB4	101062	219519	1.43%	0.106%
62	12.675	1625	1630	1635	rBV5	72800	151347	0.99%	0.073%
63	12.722	1635	1638	1647	rVB5	67538	123604	0.81%	0.060%
64	13.128	1701	1707	1713	rBV	206053	317209	2.07%	0.153%
65	13.463	1759	1764	1772	rBV2	274242	585964	3.83%	0.283%

66	13.675	1792	1800	1805	rBV	767456	1113529	7.27%	0.538%
67	13.863	1828	1832	1838	rVB	74846	112179	0.73%	0.054%
68	14.045	1856	1863	1869	rBV	7365475	10375325	67.74%	5.012%
69	14.169	1880	1884	1888	rBV2	70305	101576	0.66%	0.049%
70	14.240	1892	1896	1903	rVB4	140772	215854	1.41%	0.104%

71	14.328	1903	1911	1918	rBV	8653236	12680060	82.79%	6.125%
72	14.481	1933	1937	1944	rVB	81985	103902	0.68%	0.050%
73	14.628	1952	1962	1969	rVB	4426572	6914311	45.14%	3.340%
74	14.845	1994	1999	2006	rBV	666588	1057974	6.91%	0.511%
75	15.263	2067	2070	2077	rVB4	69116	113071	0.74%	0.055%

76	15.440	2096	2100	2108	rVB	103557	148083	0.97%	0.072%
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77	15.622	2124	2131	2137	rBV	10662364	15154530	98.94%	7.320%
78	15.687	2137	2142	2147	rVB3	245864	354600	2.32%	0.171%
79	15.745	2147	2152	2159	rBV	3507915	4799248	31.33%	2.318%
80	15.822	2161	2165	2168	rVB	299057	341194	2.23%	0.165%
81	16.151	2217	2221	2226	rBV2	107198	156994	1.03%	0.076%
82	16.486	2276	2278	2285	rVB3	64690	116524	0.76%	0.056%
83	16.551	2286	2289	2294	rBV2	47101	83083	0.54%	0.040%
84	16.822	2332	2335	2340	rBV	75030	108151	0.71%	0.052%
85	17.392	2426	2432	2437	rBV	5122146	6932499	45.26%	3.349%
86	17.434	2437	2439	2443	rVB	108861	124730	0.81%	0.060%
87	17.492	2443	2449	2458	rBV2	10945232	15288762	99.82%	7.385%
88	17.751	2489	2493	2497	rBV	217161	269315	1.76%	0.130%
89	18.233	2571	2575	2582	rBV2	635698	956239	6.24%	0.462%
90	19.051	2710	2714	2721	rVB2	231253	359598	2.35%	0.174%
91	19.298	2753	2756	2760	rBV	119698	150730	0.98%	0.073%
92	19.369	2765	2768	2772	rVB	127164	151217	0.99%	0.073%
93	19.469	2782	2785	2796	rBV2	377345	521066	3.40%	0.252%
94	19.733	2824	2830	2835	rBV	11602318	15316207	100.00%	7.398%
95	20.233	2912	2915	2920	rBV	265933	320484	2.09%	0.155%
96	20.927	3030	3033	3037	rBV	177559	209489	1.37%	0.101%
97	21.157	3069	3072	3080	rVB3	263119	370999	2.42%	0.179%
98	21.486	3123	3128	3134	rBV	4194689	5284037	34.50%	2.552%
99	23.857	3523	3531	3547	rVB2	5978251	12087636	78.92%	5.839%
100	24.021	3552	3559	3570	rVB	2392788	4933192	32.21%	2.383%

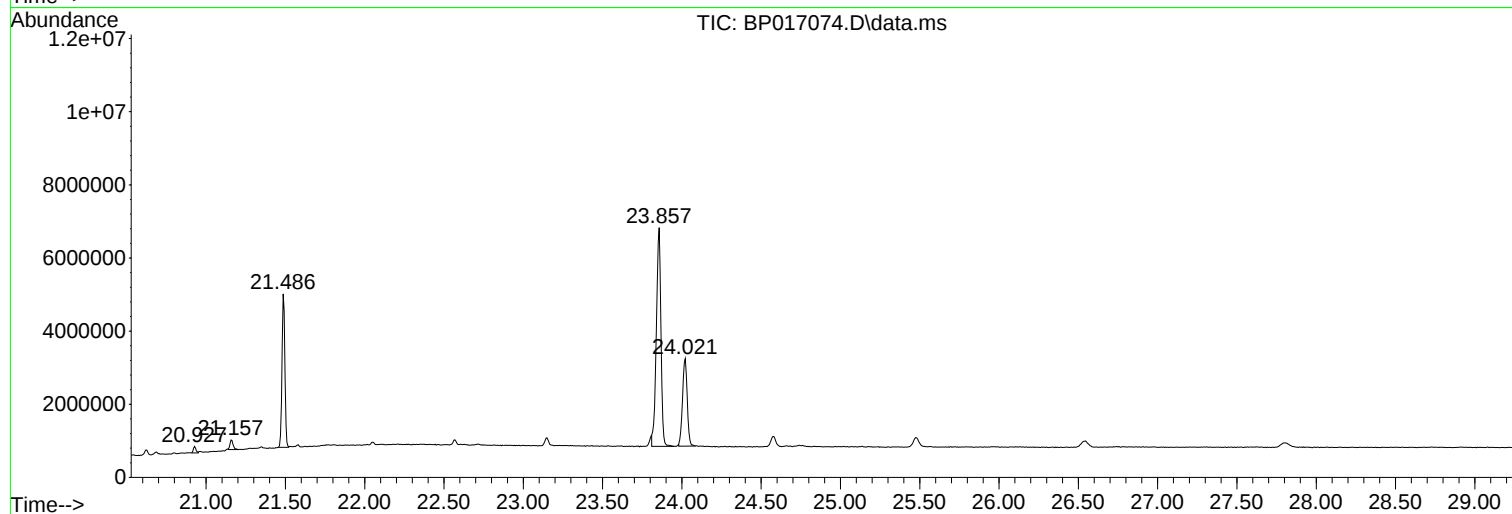
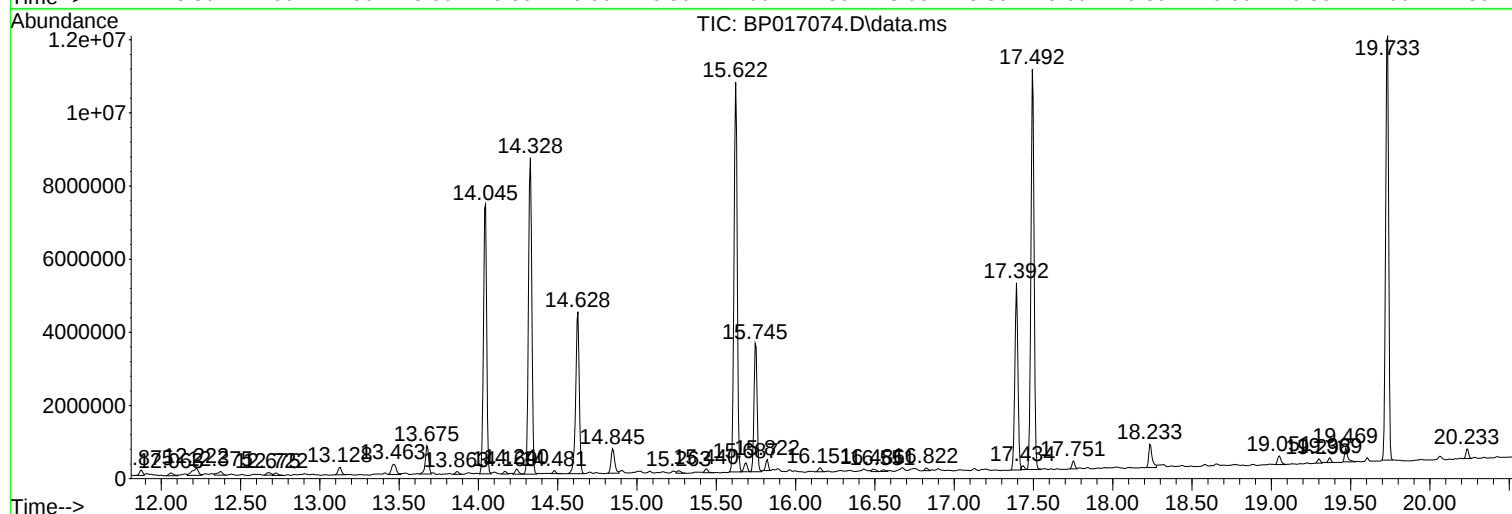
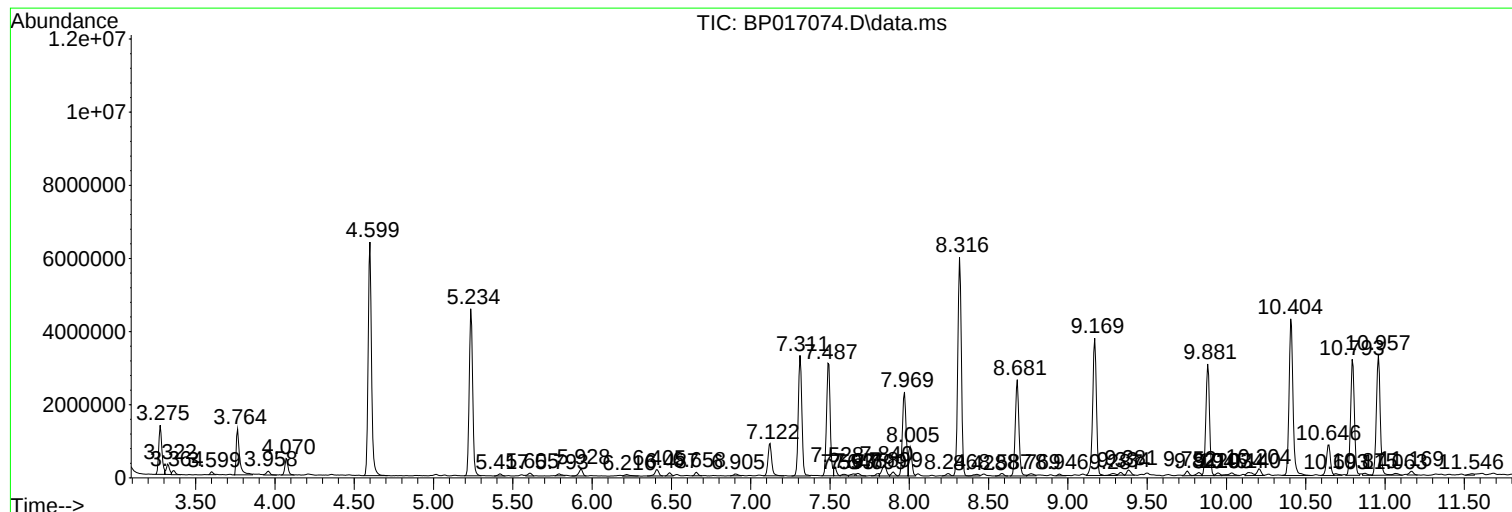
Sum of corrected areas: 207018796

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

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



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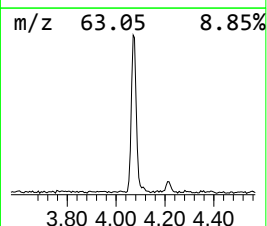
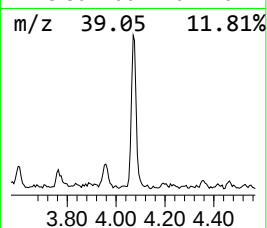
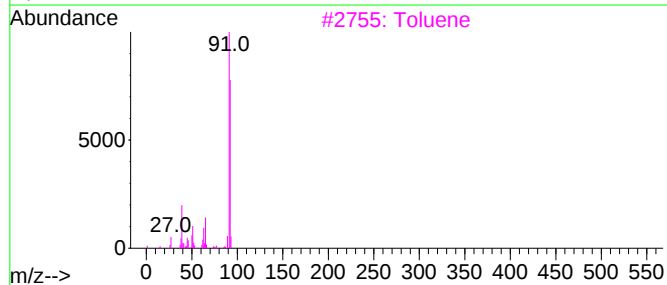
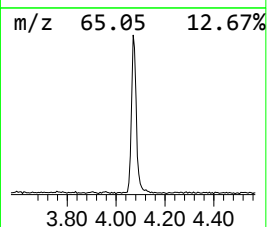
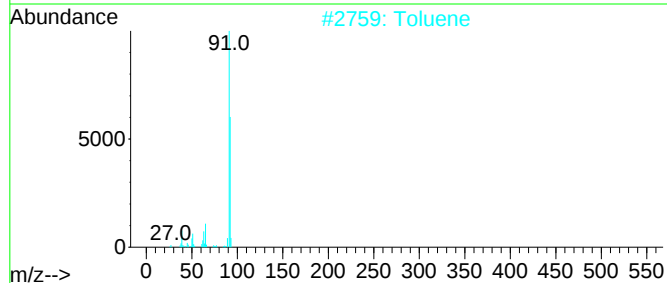
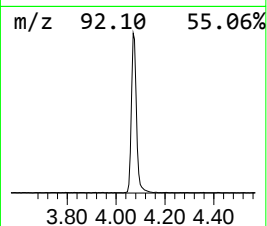
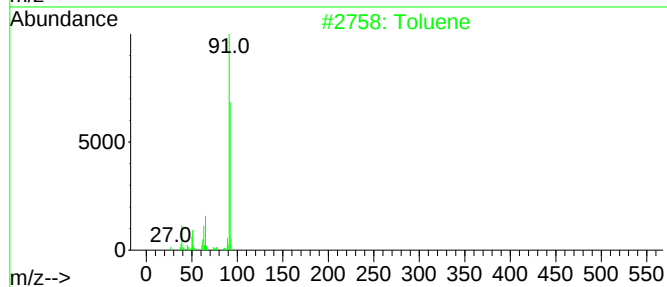
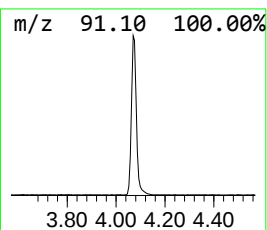
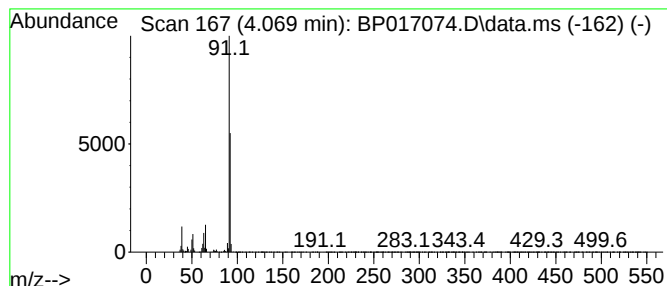
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Toluene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.069	3.32 ng/ul	628153	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene			92	C7H8	000108-88-3	95
2	Toluene			92	C7H8	000108-88-3	93
3	Toluene			92	C7H8	000108-88-3	90
4	Toluene			92	C7H8	000108-88-3	90
5	1,3,5-Cycloheptatriene			92	C7H8	000544-25-2	90



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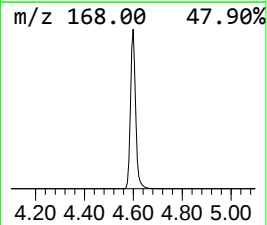
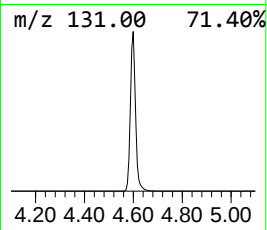
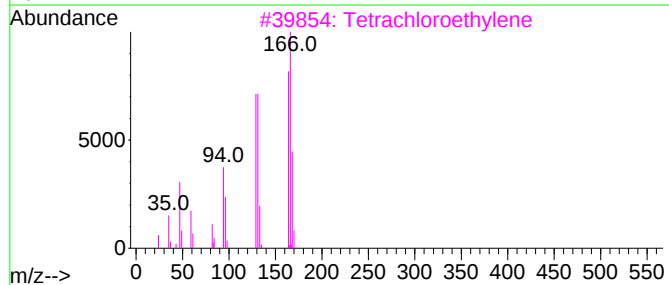
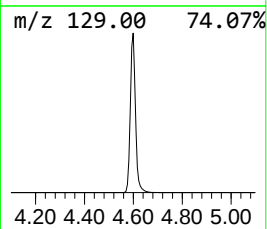
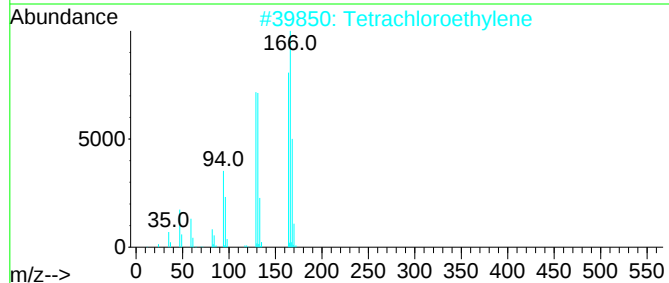
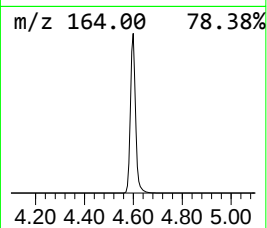
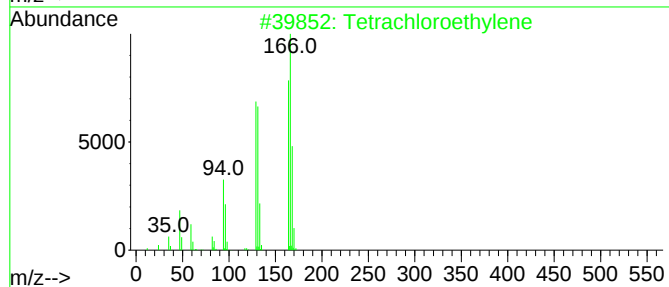
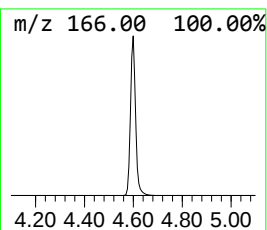
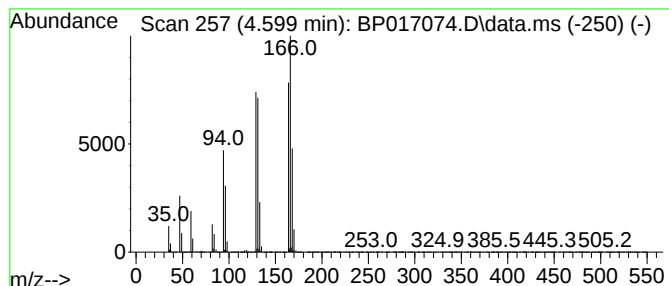
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.599	49.77 ng/ul	9411730	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	35



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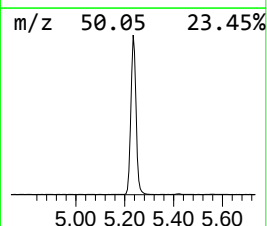
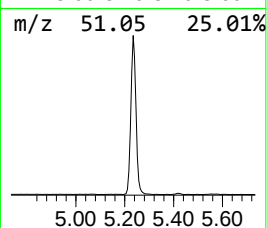
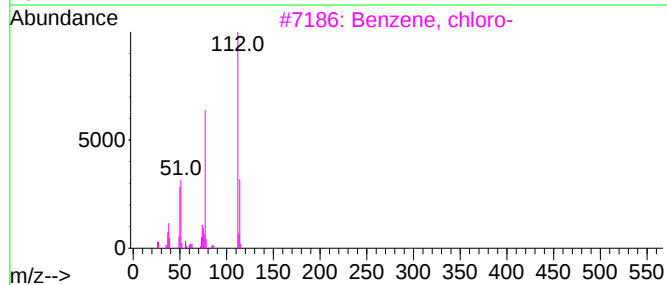
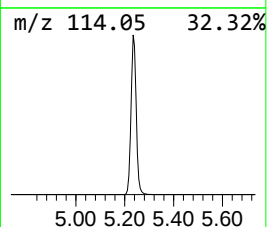
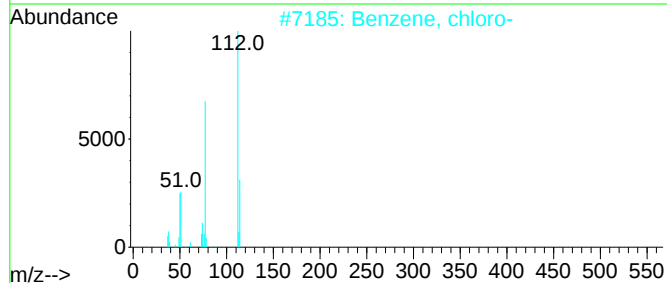
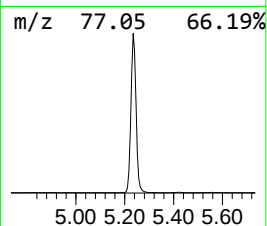
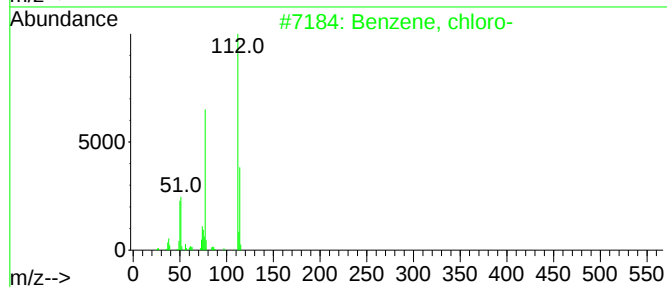
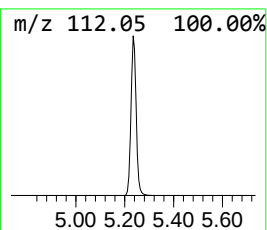
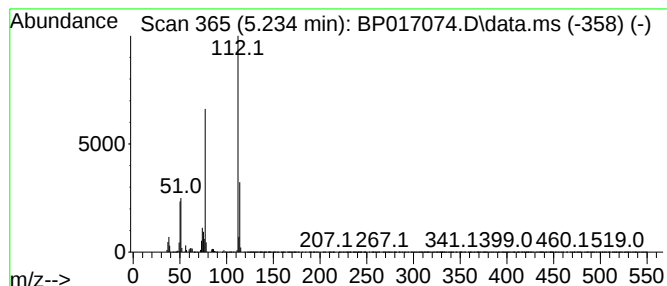
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.234	35.56 ng/ul	6724070	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	96
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
Operator : MA/JU
Sample : 04227-11
Misc :
ALS Vial : 86 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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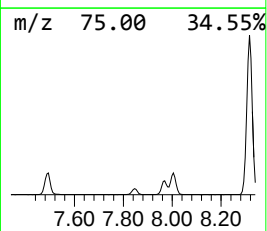
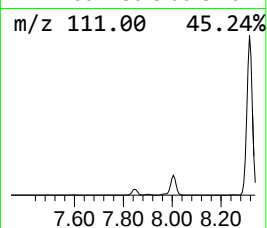
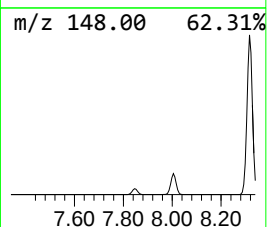
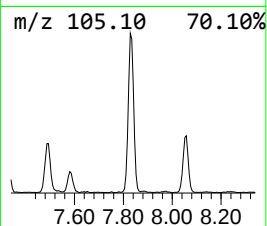
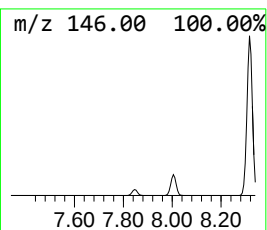
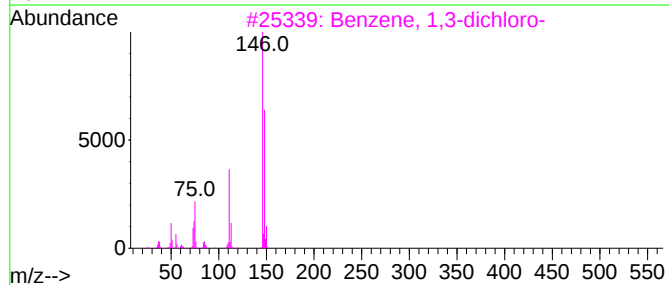
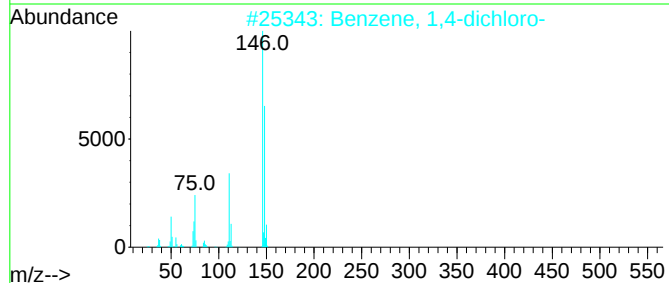
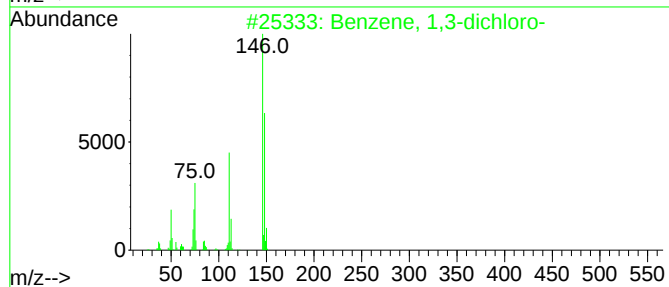
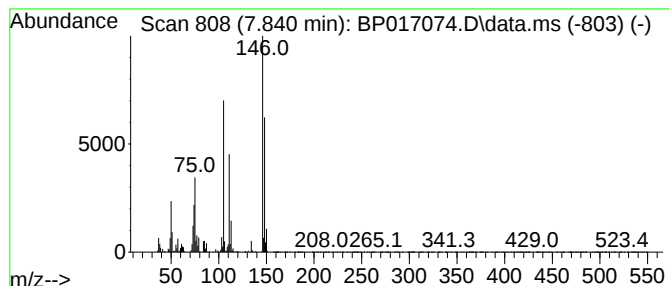
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.840	2.97 ng/ul	562534	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
Operator : MA/JU
Sample : 04227-11
Misc :
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Instrument :
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ClientSampleId :
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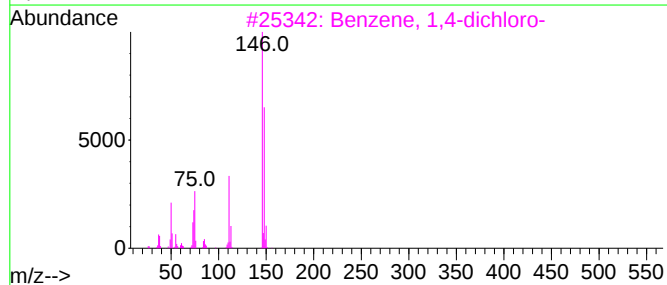
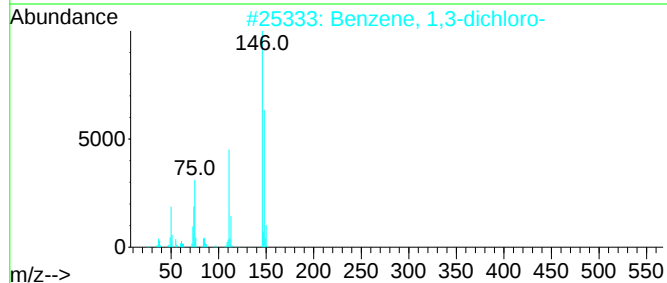
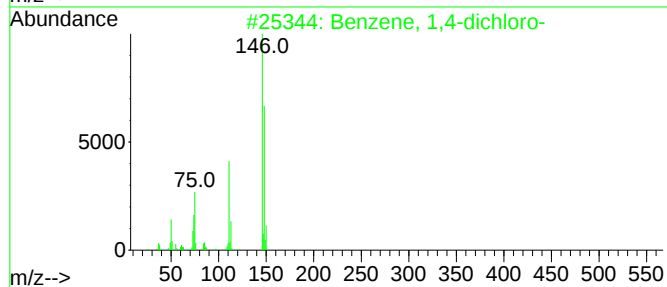
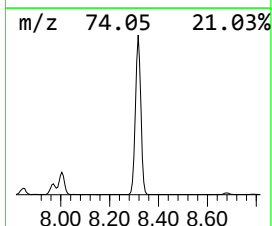
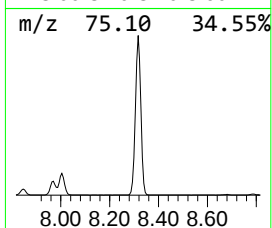
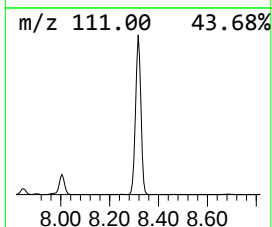
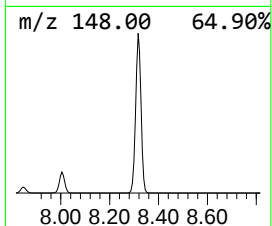
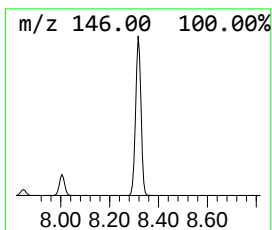
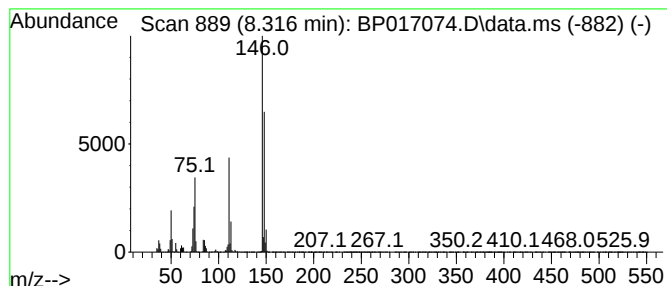
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,4-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.316	49.15 ng/ul	9293320	1,4-Dichlorobenzene-d4	7.969

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
Operator : MA/JU
Sample : 04227-11
Misc :
ALS Vial : 86 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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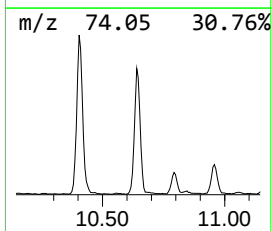
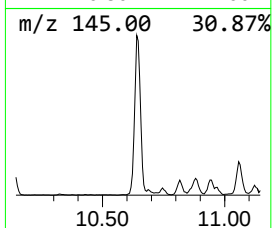
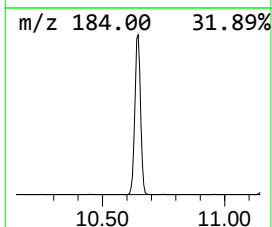
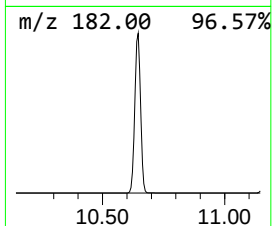
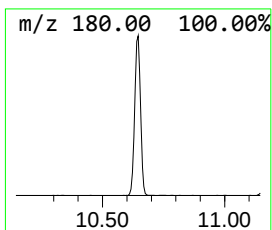
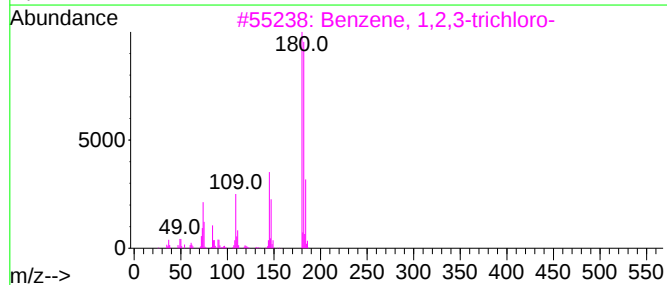
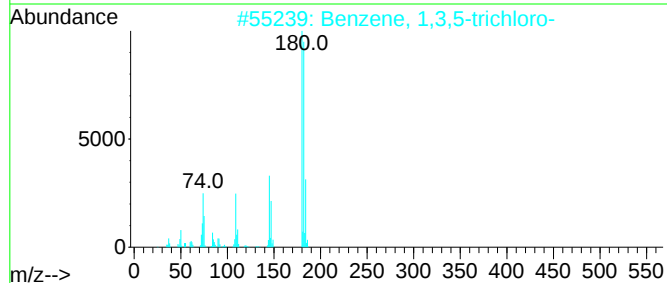
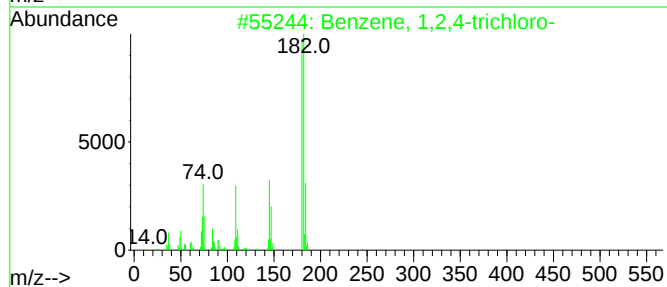
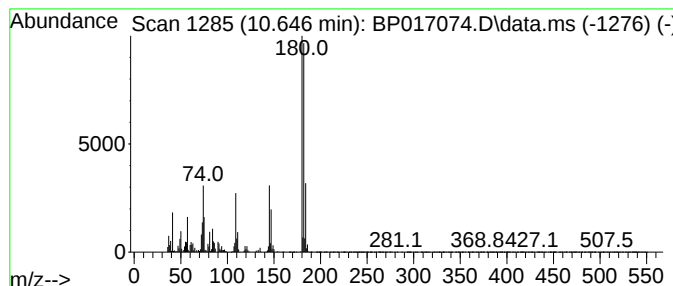
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,4-trichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.646	5.43 ng/ul	1432530	Naphthalene-d8	10.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
Operator : MA/JU
Sample : 04227-11
Misc :
ALS Vial : 86 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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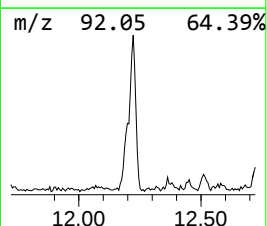
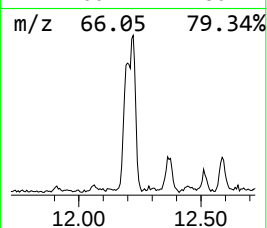
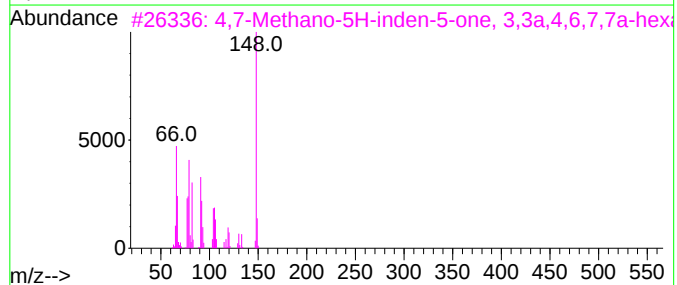
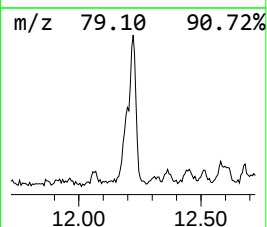
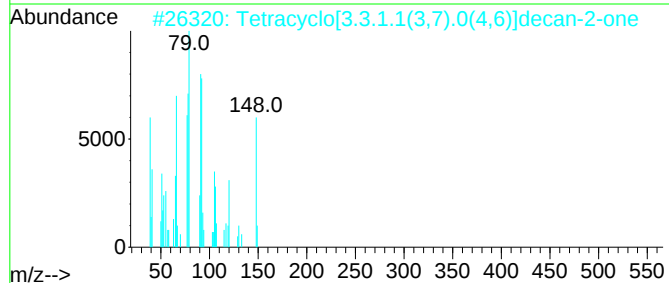
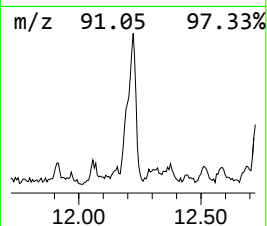
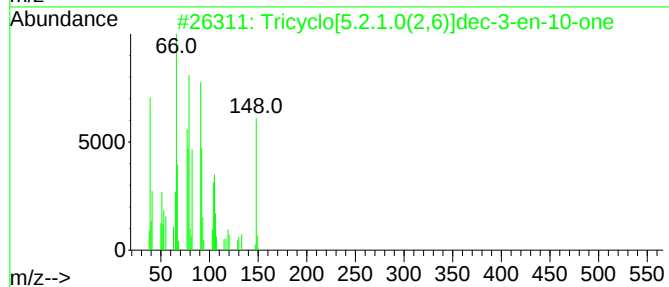
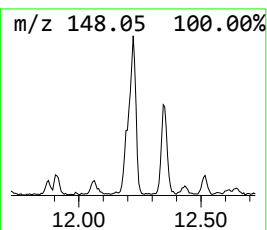
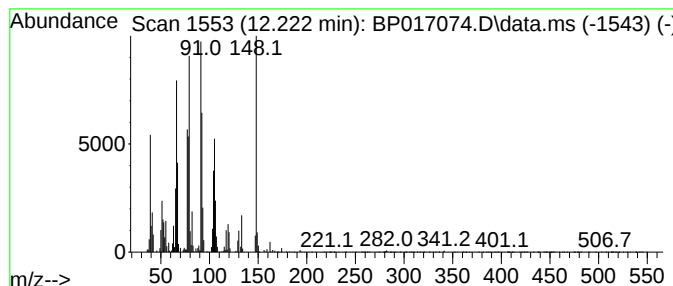
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Tricyclo[5.2.1.0(2,6)]dec-3... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.222	2.03 ng/ul	534616	Naphthalene-d8	10.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricyclo[5.2.1.0(2,6)]dec-3-en-1...	148	C10H12O	1000289-10-5	95
2			Tetracyclo[3.3.1.1(3,7).0(4,6)]d...	148	C10H12O	052750-53-5	93
3			4,7-Methano-5H-inden-5-one, 3,3a...	148	C10H12O	014888-58-5	91
4			4,7-Methano-5H-inden-5-one, 3,3a...	148	C10H12O	014888-58-5	81
5			1H-Indene, 3a,4,7,7a-tetrahydro-...	120	C9H12	066563-20-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
Operator : MA/JU
Sample : 04227-11
Misc :
ALS Vial : 86 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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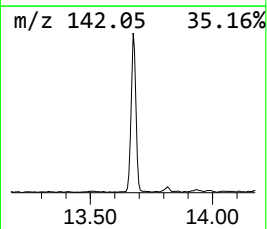
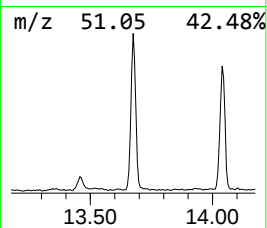
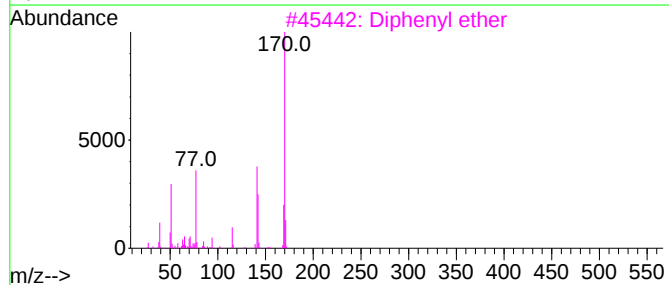
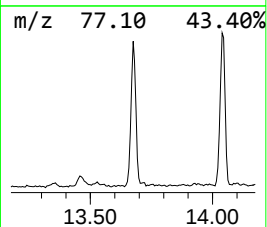
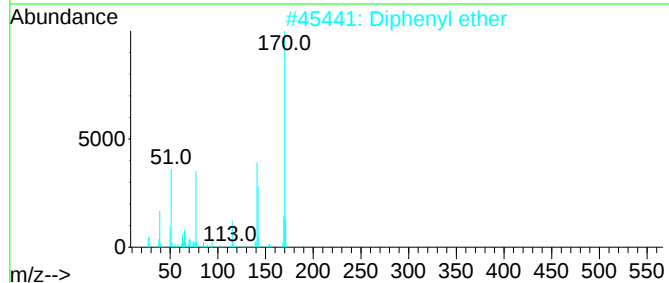
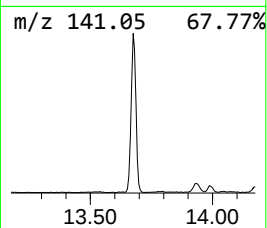
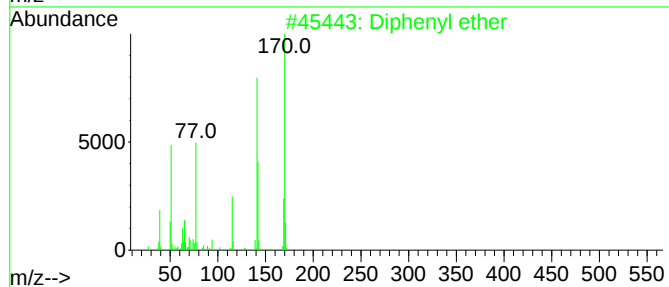
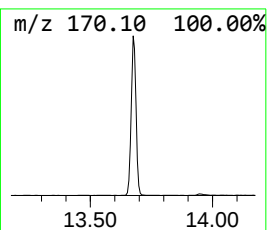
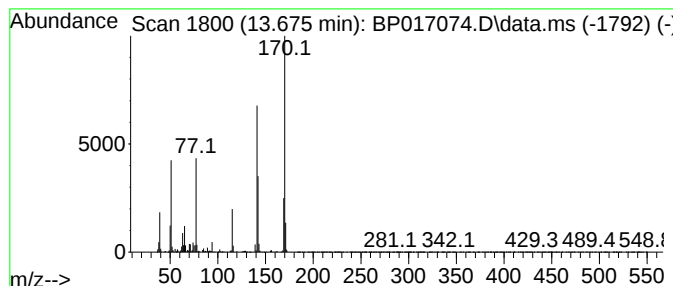
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Diphenyl ether Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.675	3.22 ng/ul	1113530	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diphenyl ether	170	C12H10O	000101-84-8	98
2			Diphenyl ether	170	C12H10O	000101-84-8	87
3			Diphenyl ether	170	C12H10O	000101-84-8	87
4			Diphenyl ether	170	C12H10O	000101-84-8	81
5			Diphenyl ether	170	C12H10O	000101-84-8	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
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Sample : 04227-11
Misc :
ALS Vial : 86 Sample Multiplier: 1

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ClientSampleId :
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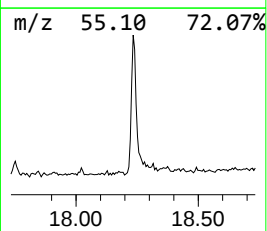
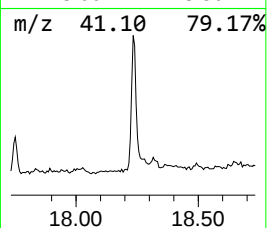
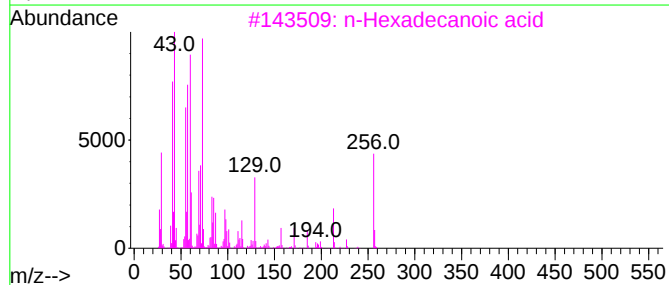
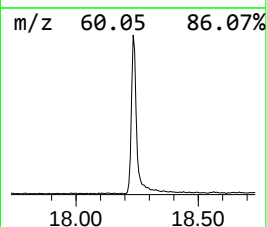
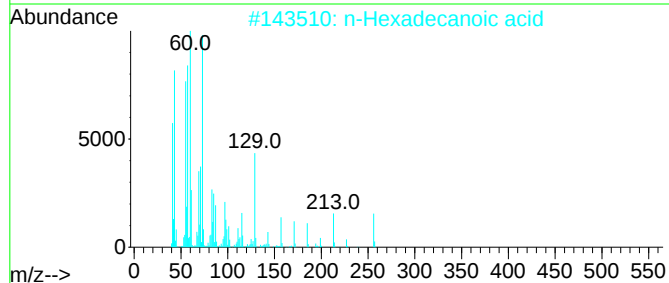
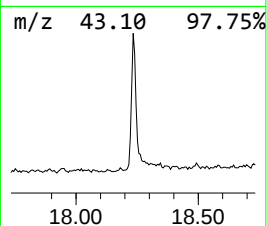
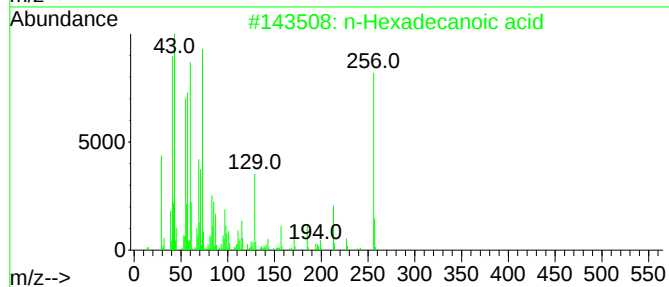
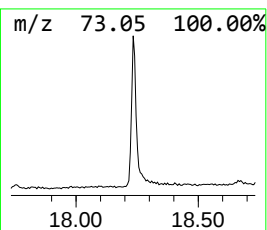
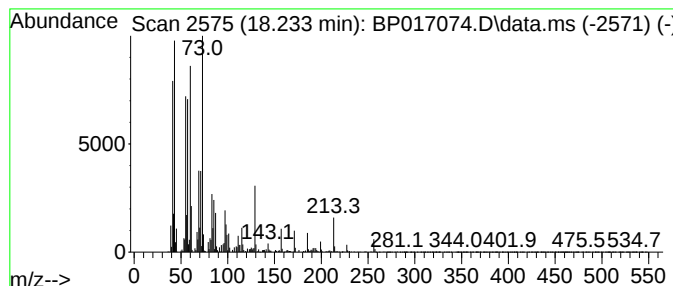
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	2.76 ng/ul	956239	Phenanthrene-d10	17.392

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	93
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090823\
Data File : BP017074.D
Acq On : 10 Sep 2023 04:23
Operator : MA/JU
Sample : 04227-11
Misc :
ALS Vial : 86 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP090523.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
Toluene	4.069	3.3	ng/ul	628153	1	7.969	3781870	20.0
Tetrachloroethy...	4.599	49.8	ng/ul	9411730	1	7.969	3781870	20.0
Benzene, chloro-	5.234	35.6	ng/ul	6724070	1	7.969	3781870	20.0
Benzene, 1,3-di...	7.840	3.0	ng/ul	562534	1	7.969	3781870	20.0
Benzene, 1,4-di...	8.316	49.1	ng/ul	9293320	1	7.969	3781870	20.0
Benzene, 1,2,4-...	10.646	5.4	ng/ul	1432530	2	10.793	5271490	20.0
Tricyclo[5.2.1....	12.222	2.0	ng/ul	534616	2	10.793	5271490	20.0
Diphenyl ether	13.675	3.2	ng/ul	1113530	3	14.628	6914310	20.0
n-Hexadecanoic ...	18.233	2.8	ng/ul	956239	4	17.392	6932500	20.0