

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017146.D
 Acq On : 13 Sep 2023 13:23
 Operator : MA/JU
 Sample : 04324-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA56

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-BP091223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017146.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	25	30	36	rBV	629634	990111	7.97%	1.347%
2	3.352	41	45	55	rVB	174491	257926	2.08%	0.351%
3	3.587	81	85	90	rBV	26147	34094	0.27%	0.046%
4	3.752	110	113	127	rBV	230988	382935	3.08%	0.521%
5	3.852	127	130	135	rVB	16872	24937	0.20%	0.034%
6	3.946	141	146	151	rVB2	33167	52104	0.42%	0.071%
7	4.058	161	165	170	rBV2	11204	16677	0.13%	0.023%
8	4.328	207	211	216	rVB	27229	36686	0.30%	0.050%
9	4.381	216	220	230	rVB2	16898	26534	0.21%	0.036%
10	4.587	245	255	261	rBV	62158	105748	0.85%	0.144%
11	5.058	329	335	339	rVV2	16576	24734	0.20%	0.034%
12	5.116	339	345	353	rVB6	16826	34408	0.28%	0.047%
13	5.393	389	392	398	rVB5	8376	12794	0.10%	0.017%
14	6.022	493	499	502	rBV2	16962	29084	0.23%	0.040%
15	6.222	527	533	540	rVB8	7593	16026	0.13%	0.022%
16	6.634	598	603	611	rBV	714407	1017144	8.19%	1.384%
17	7.105	677	683	685	rBV	81697	121402	0.98%	0.165%
18	7.140	685	689	698	rVV	138361	261630	2.11%	0.356%
19	7.293	707	715	725	rBV	714607	1111517	8.95%	1.512%
20	7.475	739	746	757	rBV	711549	1118726	9.01%	1.522%
21	7.952	820	827	836	rBV	615057	970827	7.82%	1.321%
22	8.693	941	953	967	rBV2	369128	1131316	9.11%	1.539%
23	9.034	1005	1011	1013	rBV5	13859	25609	0.21%	0.035%
24	9.069	1013	1017	1023	rVB2	24814	44407	0.36%	0.060%
25	9.151	1023	1031	1040	rBV	811338	1303674	10.50%	1.774%
26	9.246	1042	1047	1050	rBV3	13865	22709	0.18%	0.031%
27	9.369	1064	1068	1072	rVB6	9047	13129	0.11%	0.018%
28	9.457	1074	1083	1094	rBV	7653373	12416085	100.00%	16.894%
29	9.704	1118	1125	1131	rVB	9922	18451	0.15%	0.025%
30	9.763	1131	1135	1138	rBV4	15473	24687	0.20%	0.034%
31	9.804	1138	1142	1146	rVV3	23457	41664	0.34%	0.057%
32	9.863	1146	1152	1164	rVB	640198	1047708	8.44%	1.426%
33	10.051	1176	1184	1189	rVB	6866	16867	0.14%	0.023%
34	10.398	1236	1243	1250	rBV	1079548	1747722	14.08%	2.378%
35	10.540	1264	1267	1277	rVB5	14342	29130	0.23%	0.040%
36	10.781	1299	1308	1317	rBV	853169	1429720	11.52%	1.945%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
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37	10.940	1328	1335	1344	rBV	590072	992047	7.99%	1.350%
38	11.110	1361	1364	1367	rBV5	9353	13425	0.11%	0.018%
39	11.534	1431	1436	1440	rBV7	9270	20318	0.16%	0.028%
40	11.604	1443	1448	1453	rVV6	17962	34277	0.28%	0.047%
41	11.663	1455	1458	1464	rVB8	12811	23375	0.19%	0.032%
42	11.798	1472	1481	1489	rBV6	43269	118058	0.95%	0.161%
43	11.945	1500	1506	1513	rVB9	12573	29309	0.24%	0.040%
44	12.051	1516	1524	1532	rVB	129970	221786	1.79%	0.302%
45	12.145	1534	1540	1545	rVB10	7343	14581	0.12%	0.020%
46	12.240	1547	1556	1561	rBV3	27767	73750	0.59%	0.100%
47	12.298	1562	1566	1570	rBV3	15906	24089	0.19%	0.033%
48	12.363	1572	1577	1579	rBV3	17537	27007	0.22%	0.037%
49	12.545	1604	1608	1609	rBV4	10393	13031	0.10%	0.018%
50	12.663	1625	1628	1633	rBV7	11382	22514	0.18%	0.031%
51	12.781	1644	1648	1654	rBV8	9249	18449	0.15%	0.025%
52	12.881	1658	1665	1674	rBV8	14327	49961	0.40%	0.068%
53	13.045	1689	1693	1697	rBV7	9537	13566	0.11%	0.018%
54	13.157	1706	1712	1721	rVB7	18330	41148	0.33%	0.056%
55	13.322	1735	1740	1747	rBV6	42373	92792	0.75%	0.126%
56	13.463	1759	1764	1769	rBV	86117	125483	1.01%	0.171%
57	13.604	1783	1788	1796	rBV	246987	386359	3.11%	0.526%
58	13.739	1805	1811	1817	rVB9	15757	40781	0.33%	0.055%
59	13.816	1819	1824	1827	rBV5	19481	37352	0.30%	0.051%
60	13.857	1827	1831	1834	rVV2	33284	48559	0.39%	0.066%
61	13.892	1834	1837	1843	rVB	72845	108413	0.87%	0.148%
62	14.028	1854	1860	1870	rBV	2296829	3038465	24.47%	4.134%
63	14.116	1872	1875	1883	rVB6	25490	45210	0.36%	0.062%
64	14.192	1884	1888	1890	rBV4	18350	28196	0.23%	0.038%
65	14.228	1890	1894	1901	rVB3	46060	68830	0.55%	0.094%
66	14.310	1901	1908	1918	rBV	2258355	3357606	27.04%	4.569%
67	14.392	1918	1922	1925	rBV3	19399	27841	0.22%	0.038%
68	14.510	1937	1942	1947	rBV5	35635	55039	0.44%	0.075%
69	14.610	1953	1959	1967	rVB	1296210	1920056	15.46%	2.613%
70	14.681	1968	1971	1976	rVB6	21698	31042	0.25%	0.042%
71	14.734	1976	1980	1985	rBV7	16112	35981	0.29%	0.049%
72	14.834	1992	1997	2005	rBV	211248	375248	3.02%	0.511%
73	14.975	2017	2021	2030	rVB	53127	105393	0.85%	0.143%
74	15.075	2032	2038	2040	rBV6	18754	35201	0.28%	0.048%
75	15.163	2047	2053	2059	rBV	419036	621294	5.00%	0.845%
76	15.245	2060	2067	2078	rVB3	85550	170776	1.38%	0.232%

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77	15.451	2099	2102	2109	rVB5	20321	35246	0.28%	0.048%
78	15.610	2122	2129	2137	rBV	3194199	4487745	36.14%	6.106%
79	15.733	2145	2150	2165	rBV	1127745	1606895	12.94%	2.186%
80	15.851	2165	2170	2171	rVV4	46905	63882	0.51%	0.087%
81	15.886	2171	2176	2183	rVV	503212	746976	6.02%	1.016%
82	15.957	2184	2188	2196	rVB4	37542	66208	0.53%	0.090%
83	16.186	2224	2227	2232	rVB7	18562	25674	0.21%	0.035%
84	16.245	2232	2237	2247	rBV	118803	206173	1.66%	0.281%
85	16.986	2360	2363	2366	rBV4	22145	34370	0.28%	0.047%
86	17.116	2383	2385	2389	rVB5	13081	13129	0.11%	0.018%
87	17.380	2424	2430	2437	rBV	1735725	2442844	19.67%	3.324%
88	17.486	2441	2448	2457	rBV	3908014	5542254	44.64%	7.541%
89	18.169	2561	2564	2569	rVB2	37358	46700	0.38%	0.064%
90	18.227	2569	2574	2583	rBV	481693	682210	5.49%	0.928%
91	19.292	2752	2755	2758	rBV	52178	58599	0.47%	0.080%
92	19.363	2764	2767	2771	rVB2	57577	72819	0.59%	0.099%
93	19.463	2780	2784	2792	rBV	320656	442165	3.56%	0.602%
94	19.722	2822	2828	2840	rBV	5495133	6919621	55.73%	9.415%
95	21.486	3123	3128	3134	rVB	1921898	2568123	20.68%	3.494%
96	23.845	3521	3529	3541	rVB2	3111767	6426597	51.76%	8.744%
97	24.010	3550	3557	3567	rVB	1237902	2539707	20.45%	3.456%

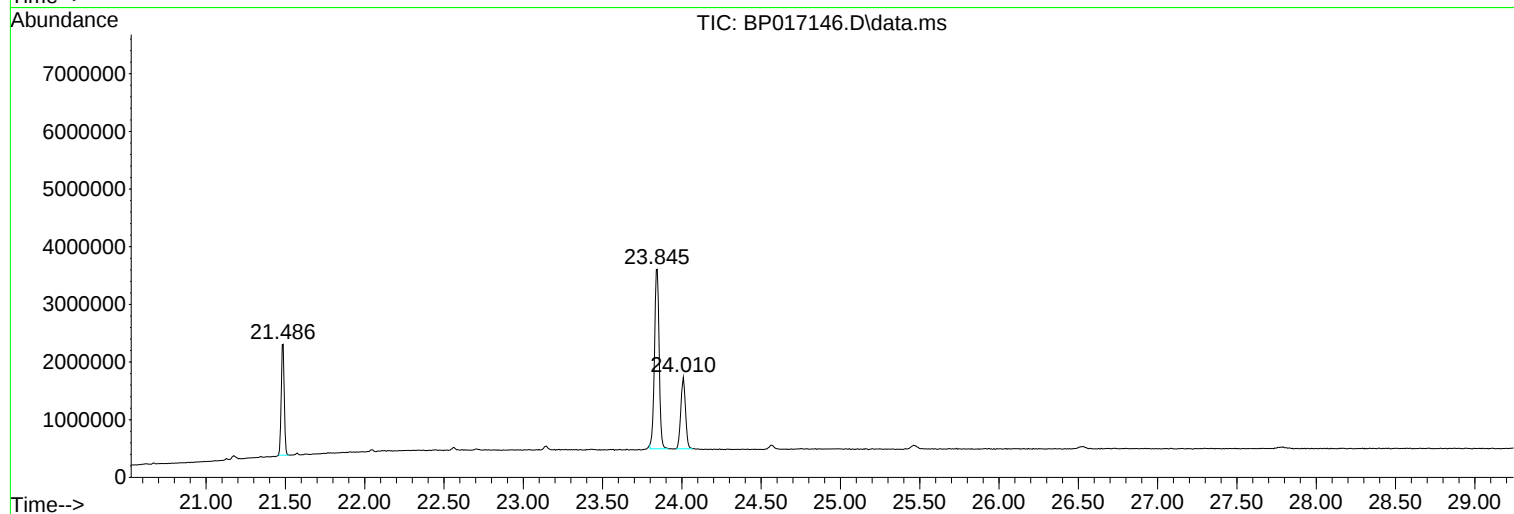
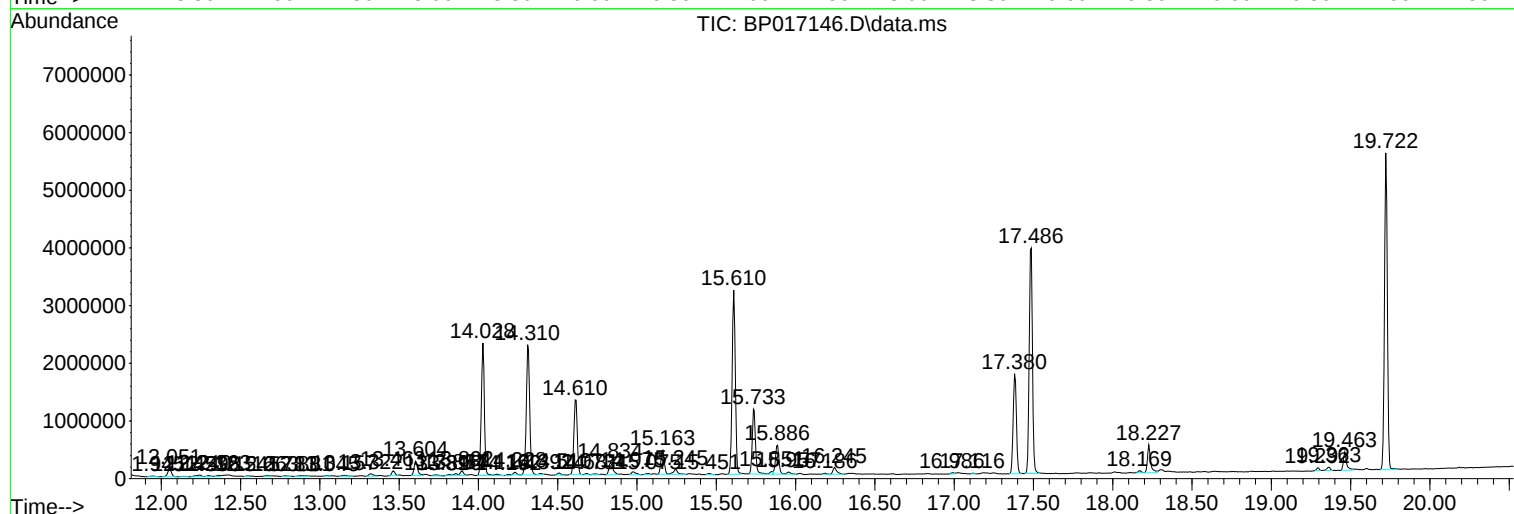
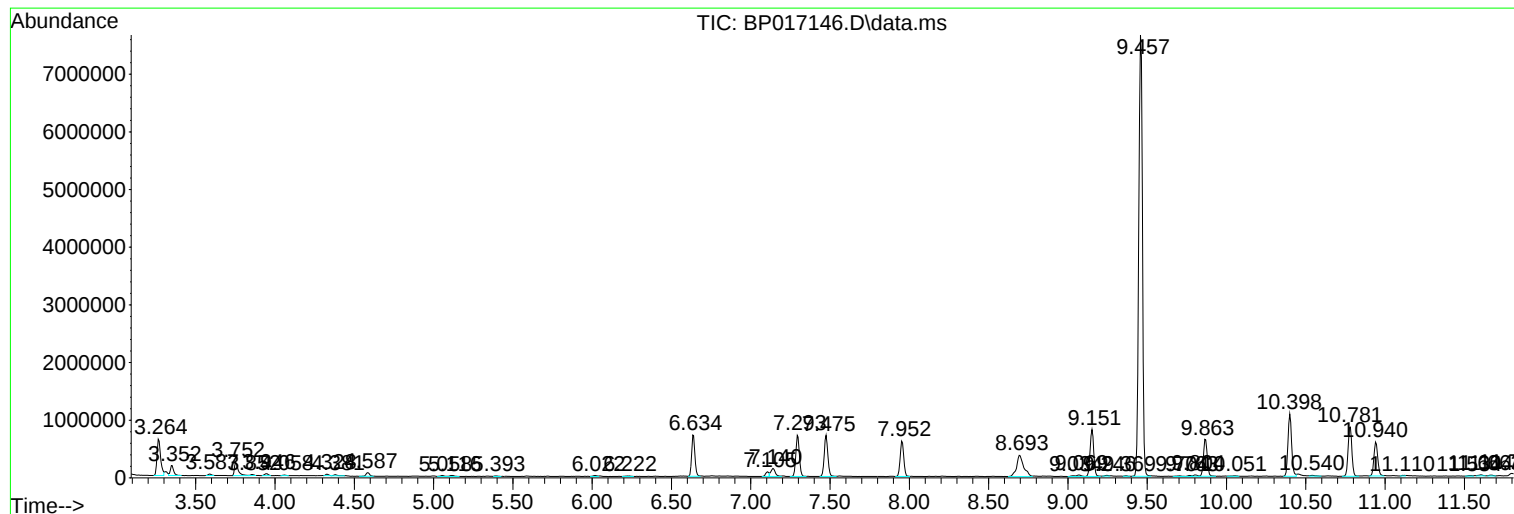
Sum of corrected areas: 73493467

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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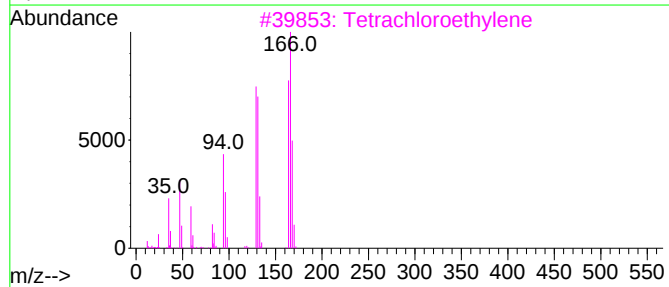
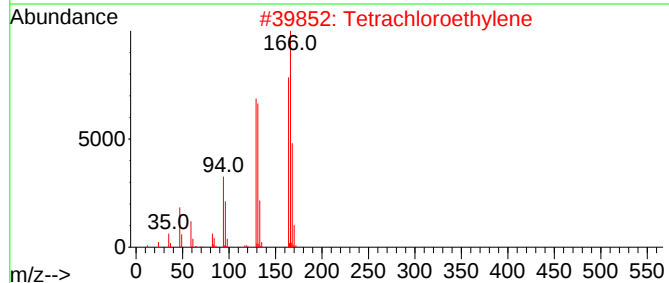
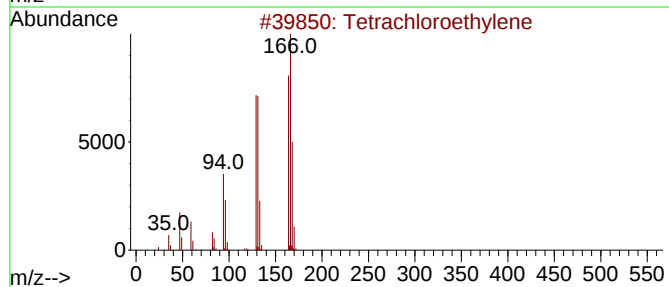
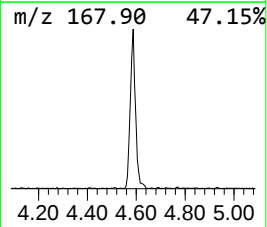
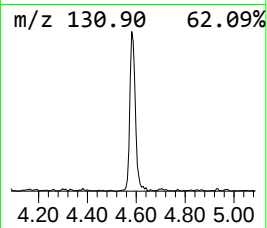
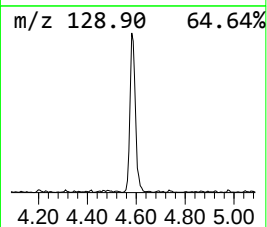
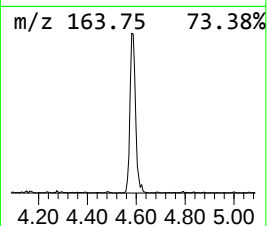
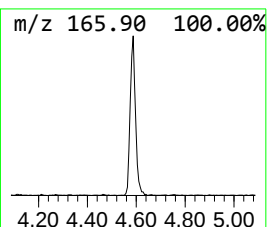
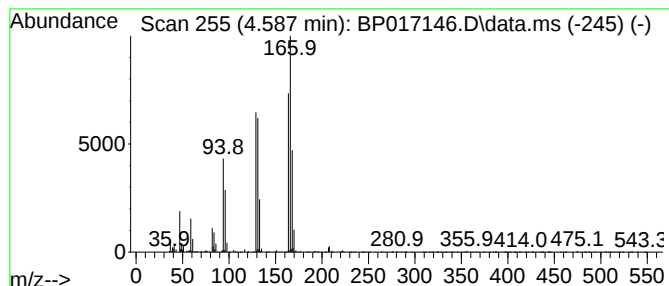
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.587	2.18 ng/ul	105748	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	95
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	32



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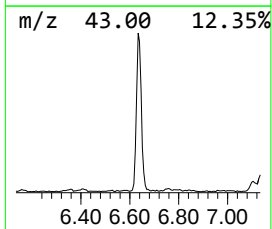
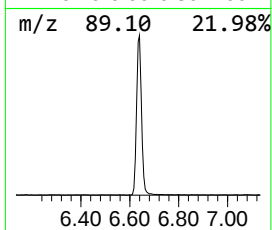
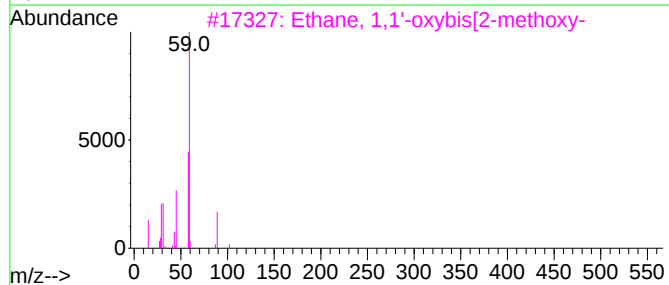
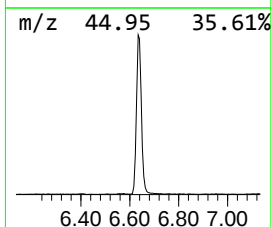
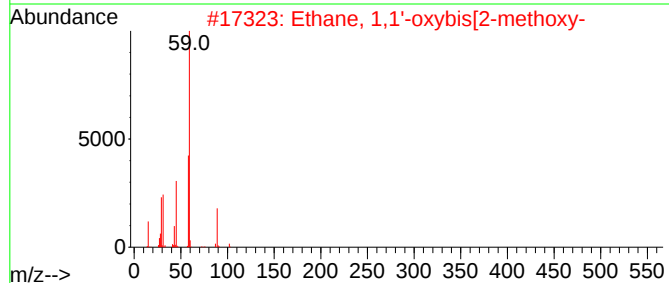
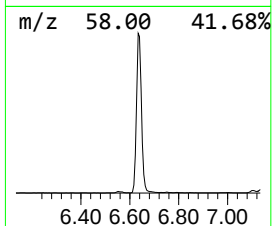
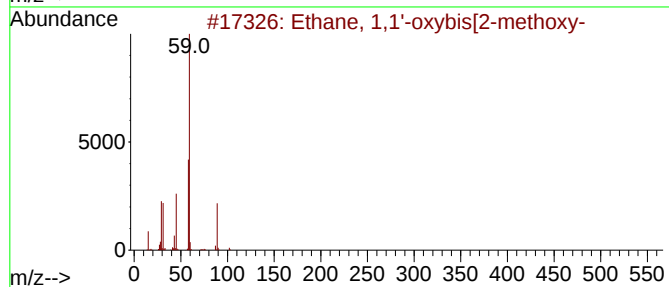
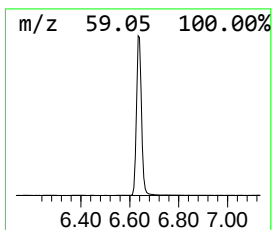
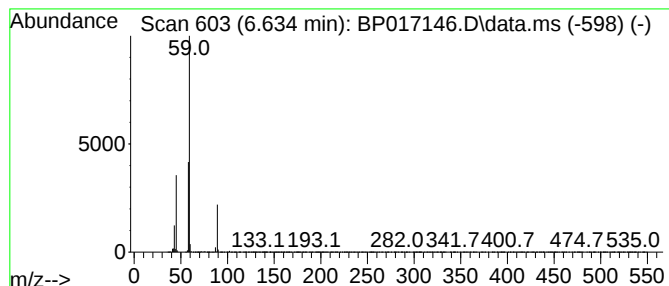
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethane, 1,1'-oxybis[2-methoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.634	20.95 ng/ul	1017140	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,1'-oxybis[2-methoxy-	134	C6H14O3	000111-96-6	83
2			Ethane, 1,1'-oxybis[2-methoxy-	134	C6H14O3	000111-96-6	83
3			Ethane, 1,1'-oxybis[2-methoxy-	134	C6H14O3	000111-96-6	83
4			Ethane, 1,1'-oxybis[2-methoxy-	134	C6H14O3	000111-96-6	64
5			Ethanol, 2-ethoxy-	90	C4H10O2	000110-80-5	9



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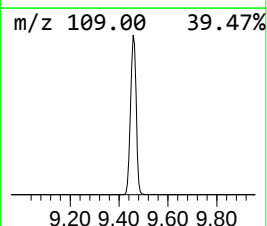
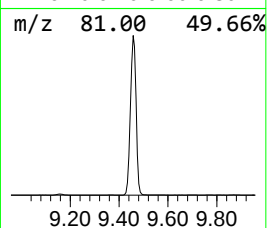
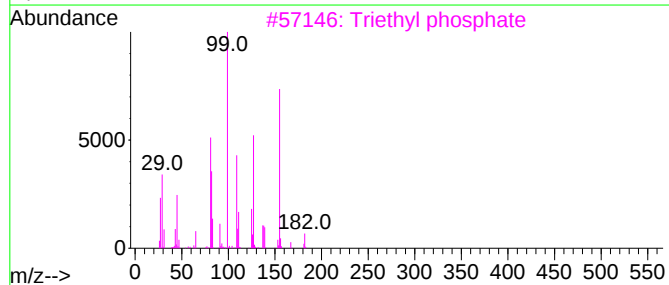
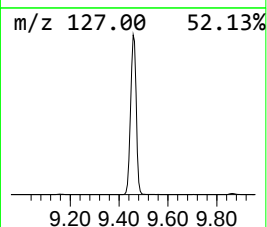
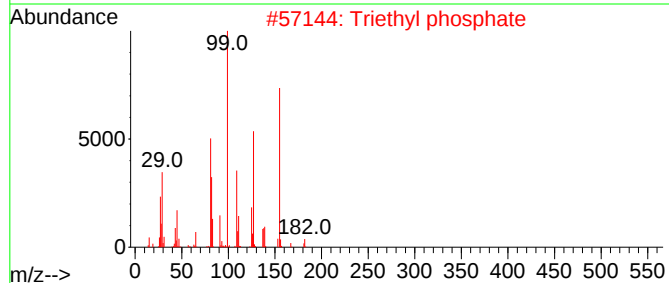
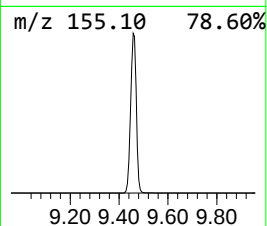
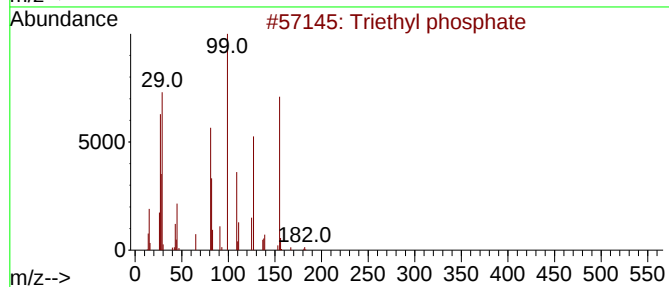
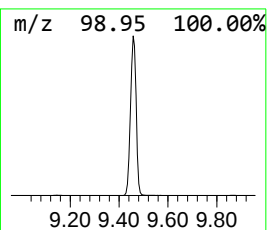
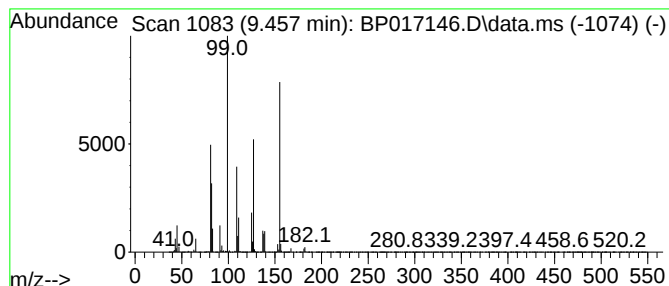
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Triethyl phosphate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.457	173.69 ng/ul	12416100	Naphthalene-d8	10.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Triethyl phosphate	182	C6H15O4P	000078-40-0	94
2			Triethyl phosphate	182	C6H15O4P	000078-40-0	94
3			Triethyl phosphate	182	C6H15O4P	000078-40-0	91
4			Triethyl phosphate	182	C6H15O4P	000078-40-0	90
5			Phosphoric acid, diethyl undecyl...	308	C15H33O4P	1000308-88-8	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
Operator : MA/JU
Sample : 04324-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA56

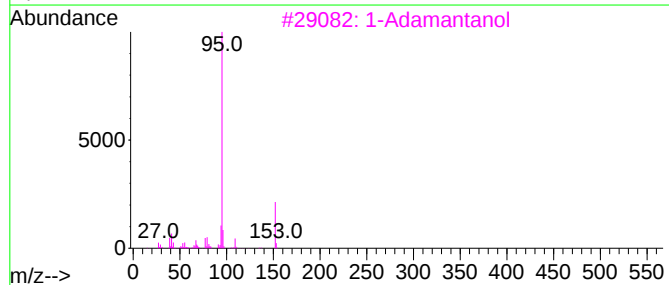
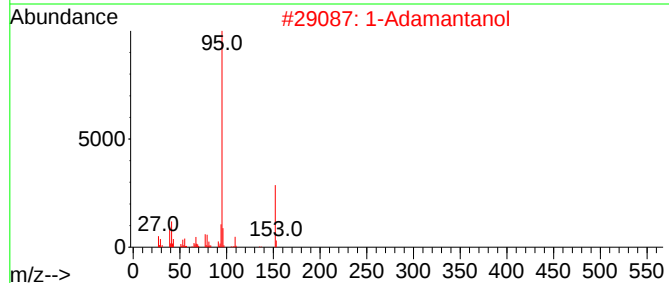
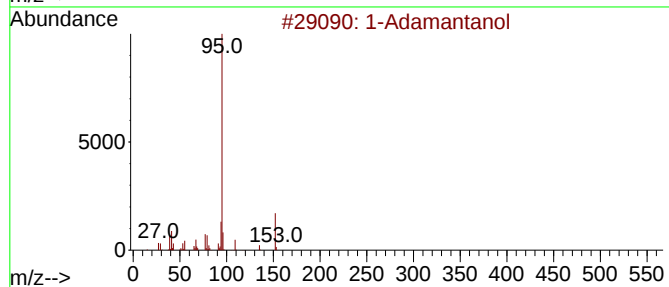
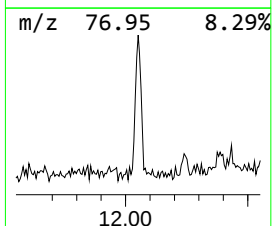
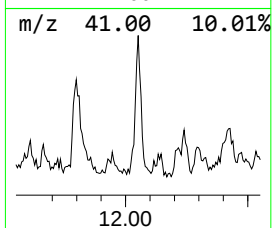
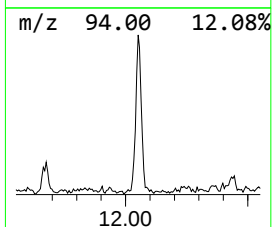
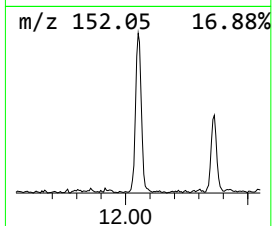
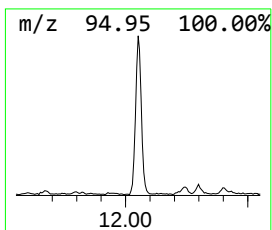
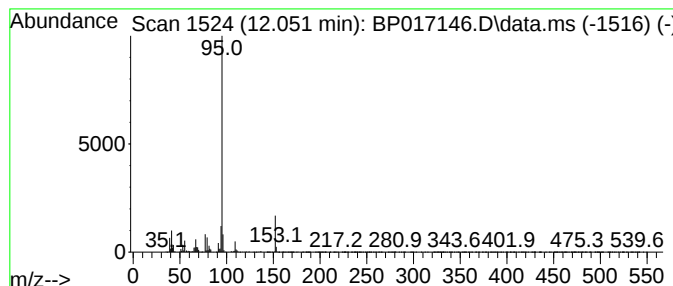
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Adamantanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.051	3.10 ng/ul	221786	Naphthalene-d8	10.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Adamantanol			152	C10H16O	000768-95-6	94
2	1-Adamantanol			152	C10H16O	000768-95-6	91
3	1-Adamantanol			152	C10H16O	000768-95-6	91
4	1-Adamantanol			152	C10H16O	000768-95-6	91
5	Succinic acid, di(hept-1,6-dien-...			306	C18H26O4	1000391-33-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
Operator : MA/JU
Sample : 04324-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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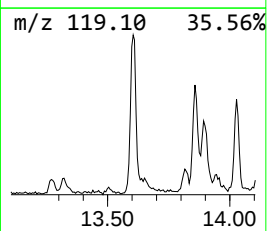
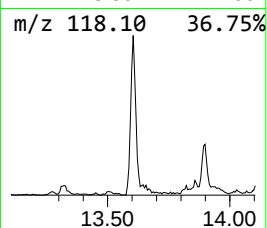
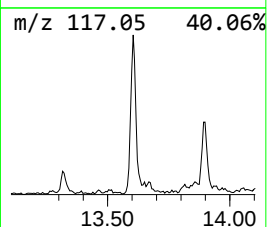
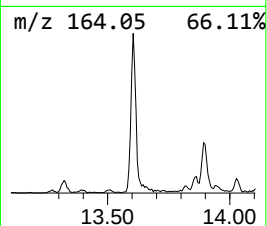
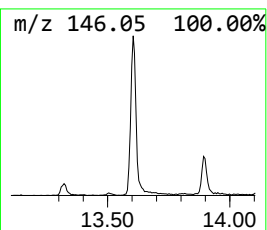
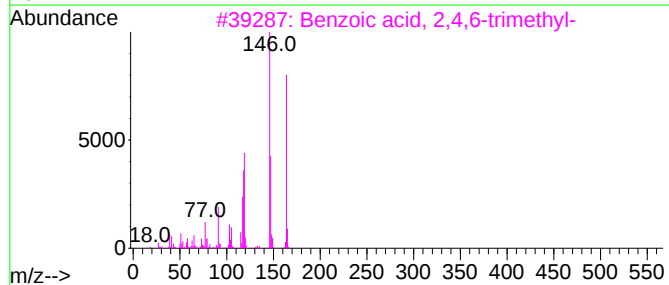
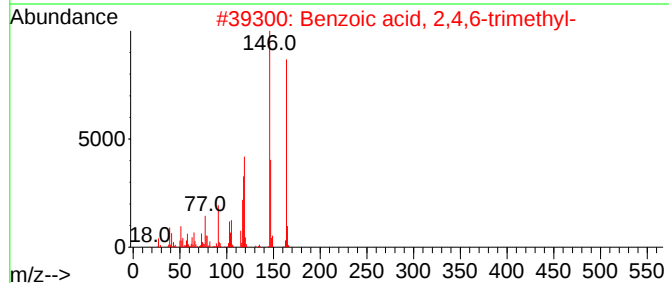
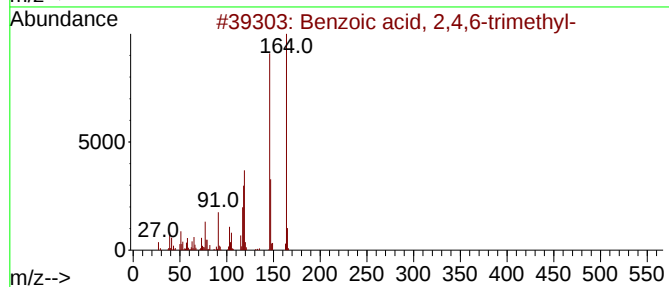
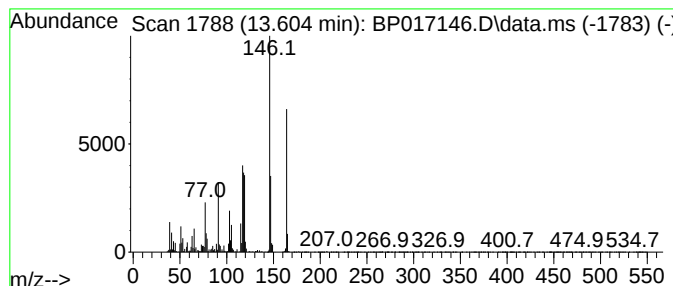
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzoic acid, 2,4,6-trimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.604	4.02 ng/ul	386359	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	81
2			Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	64
3			Benzoic acid, 2,4,6-trimethyl-	164	C10H12O2	000480-63-7	60
4			p-Coumaric acid	164	C9H8O3	007400-08-0	38
5			p-Coumaric acid, trans	164	C9H8O3	000501-98-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
Operator : MA/JU
Sample : 04324-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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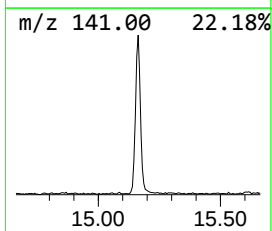
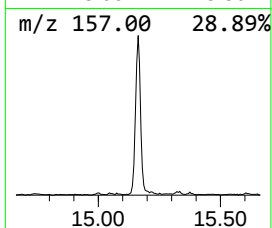
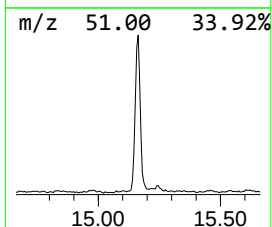
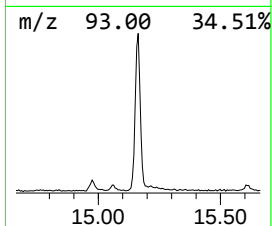
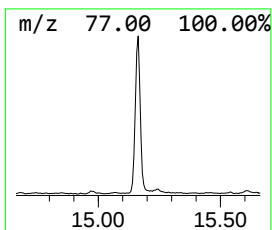
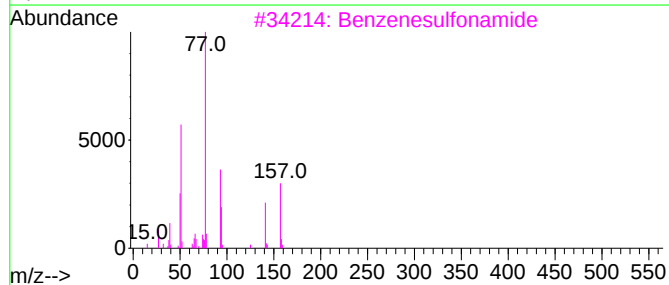
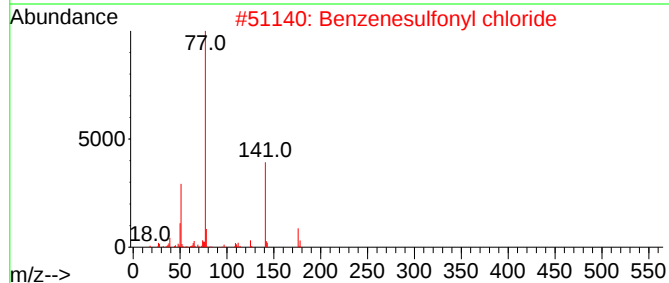
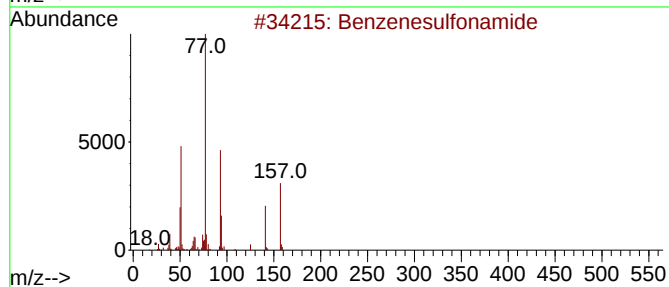
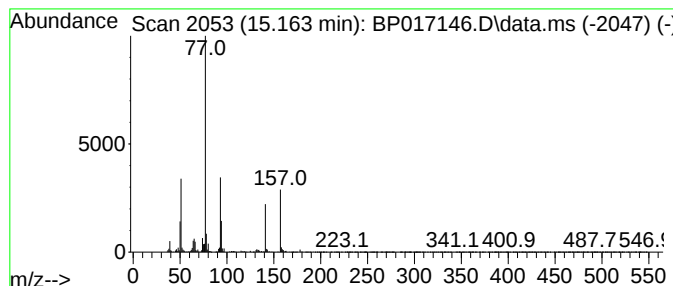
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzenesulfonamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.163	6.47 ng/ul	621294	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide	157	C6H7NO2S	000098-10-2	96
2			Benzenesulfonyl chloride	176	C6H5ClO2S	000098-09-9	50
3			Benzenesulfonamide	157	C6H7NO2S	000098-10-2	47
4			(Phenylsulfonyl)acetonitrile	181	C8H7NO2S	007605-28-9	47
5			4H-1,2,4-Triazole-4-benzenesulfo...	224	C8H8N4O2S	029982-62-5	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
Operator : MA/JU
Sample : 04324-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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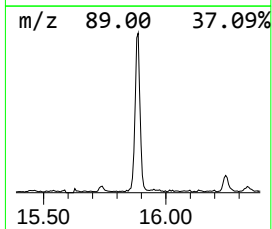
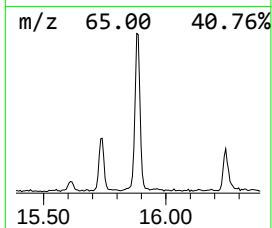
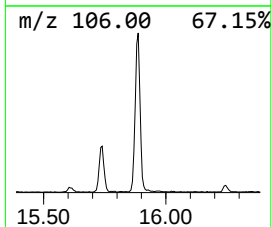
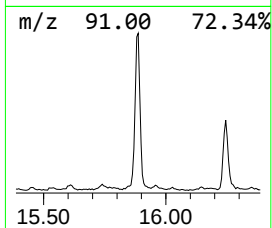
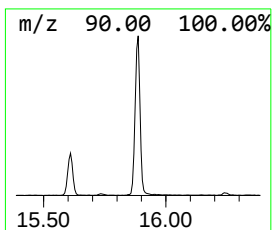
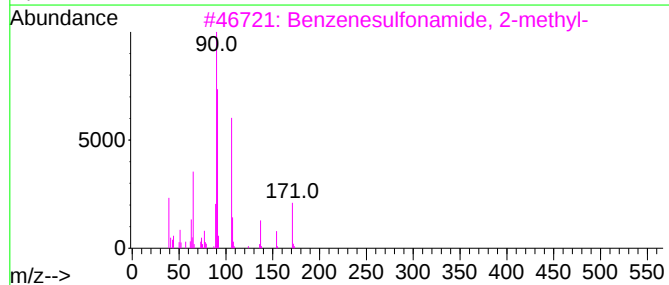
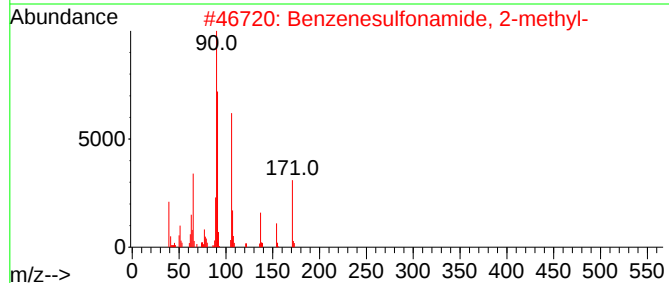
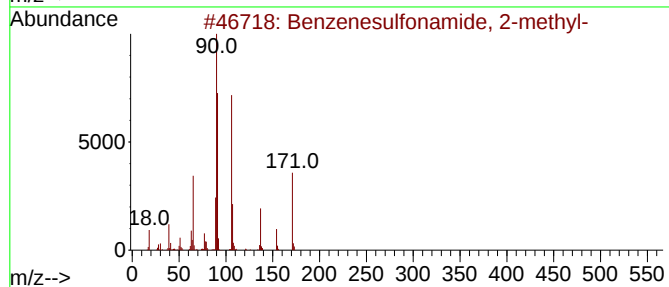
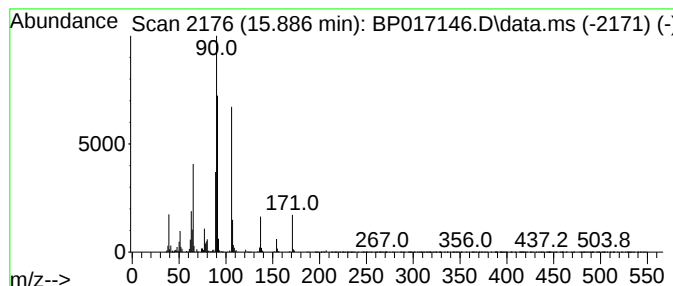
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzenesulfonamide, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.886	7.78 ng/ul	746976	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, 2-methyl-	171	C7H9NO2S	000088-19-7	96
2			Benzenesulfonamide, 2-methyl-	171	C7H9NO2S	000088-19-7	94
3			Benzenesulfonamide, 2-methyl-	171	C7H9NO2S	000088-19-7	94
4			Benzenesulfonamide, 2-methyl-	171	C7H9NO2S	000088-19-7	91
5			4-(2-Hydroxyethyl)benzenesulfona...	201	C8H11NO3S	000829-71-0	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
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Sample : 04324-06
Misc :
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Instrument :
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ClientSampleId :
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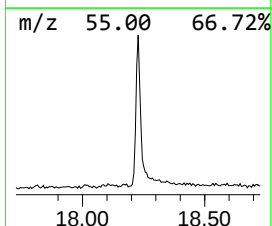
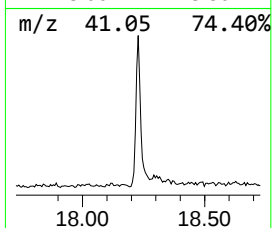
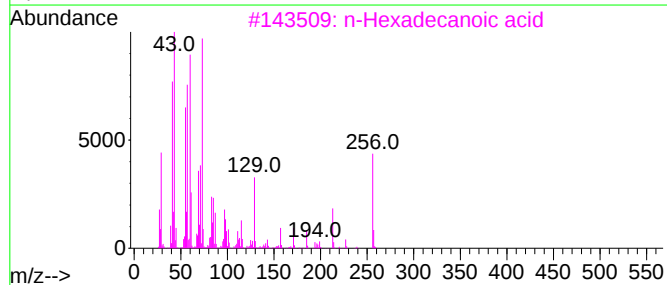
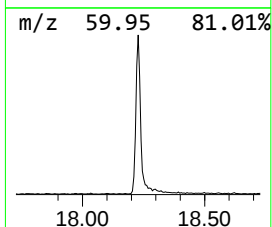
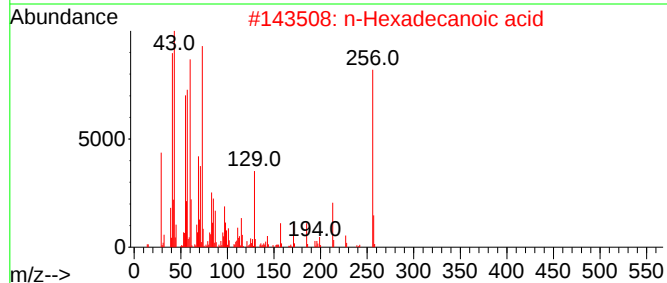
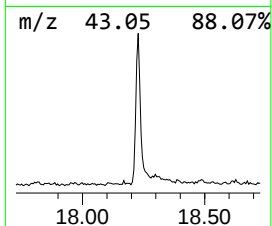
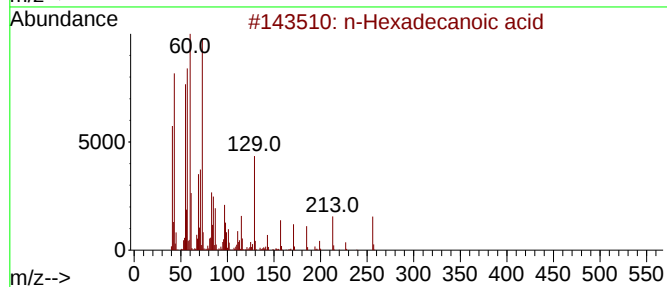
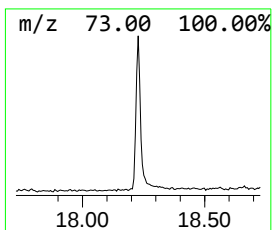
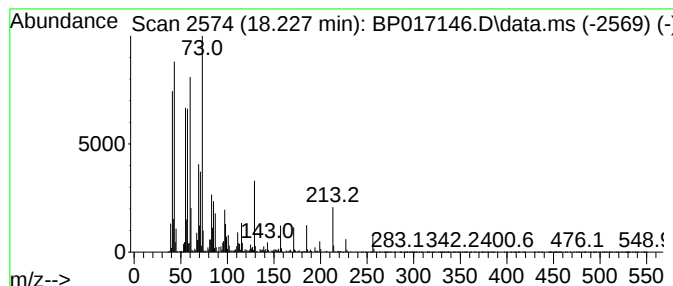
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.227	5.59 ng/ul	682210	Phenanthrene-d10	17.381

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	95
5			Tridecanoic acid	214	C13H26O2	000638-53-9	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
Operator : MA/JU
Sample : 04324-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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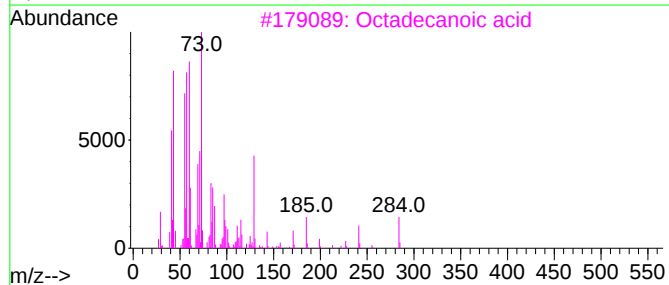
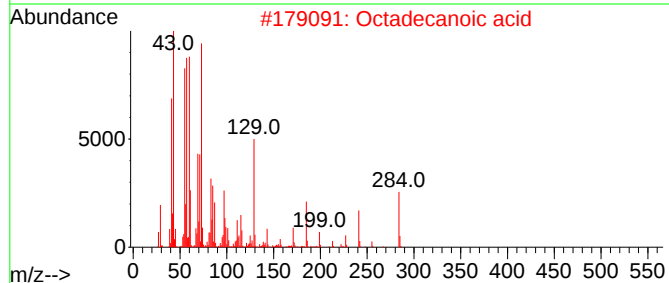
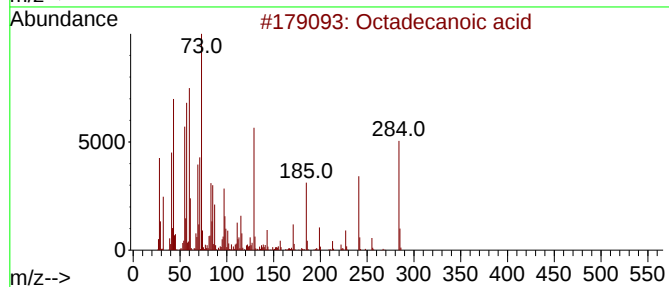
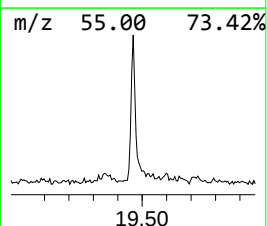
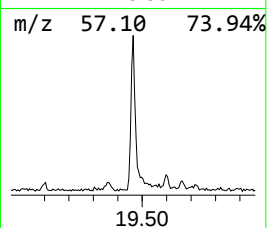
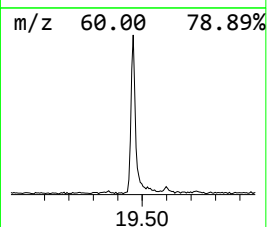
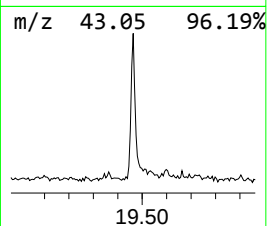
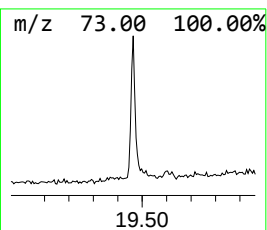
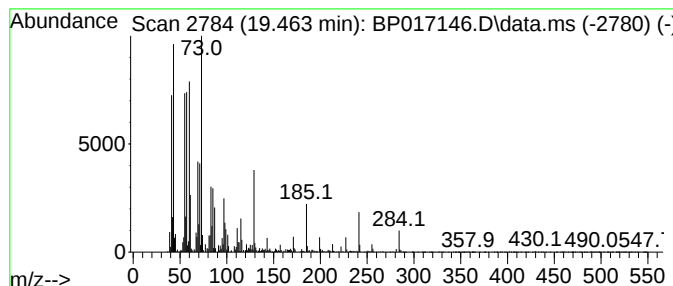
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.463	3.44 ng/ul	442165	Chrysene-d12	21.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid			284	C18H36O2	000057-11-4	99
2	Octadecanoic acid			284	C18H36O2	000057-11-4	99
3	Octadecanoic acid			284	C18H36O2	000057-11-4	98
4	Octadecanoic acid			284	C18H36O2	000057-11-4	95
5	Octadecanoic acid			284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
Data File : BP017146.D
Acq On : 13 Sep 2023 13:23
Operator : MA/JU
Sample : 04324-06
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA56

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.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
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Ethane, 1,1'-ox...	6.634	20.9	ng/ul	1017140	1	7.952	970827	20.0
Triethyl phosphate	9.457	173.7	ng/ul	12416100	2	10.781	1429720	20.0
1-Adamantanol	12.051	3.1	ng/ul	221786	2	10.781	1429720	20.0
Benzoic acid, 2...	13.604	4.0	ng/ul	386359	3	14.610	1920060	20.0
Benzenesulfonamide	15.163	6.5	ng/ul	621294	3	14.610	1920060	20.0
Benzenesulfonam...	15.886	7.8	ng/ul	746976	3	14.610	1920060	20.0
n-Hexadecanoic ...	18.227	5.6	ng/ul	682210	4	17.381	2442840	20.0
Octadecanoic acid	19.463	3.4	ng/ul	442165	5	21.486	2568120	20.0