

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
 Data File : BP020353.D
 Acq On : 13 May 2024 12:47
 Operator : MA/JU
 Sample : P2372-23DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43Y3DL

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-BP050124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020353.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.911	137	140	148	rVB5	5231	8199	0.26%	0.018%
2	4.405	219	224	225	rBV5	2419	3558	0.11%	0.008%
3	6.834	626	637	652	rBV2	199450	526508	16.88%	1.151%
4	6.975	656	661	670	rVV	30414	63433	2.03%	0.139%
5	7.063	670	676	686	rVV	53023	106097	3.40%	0.232%
6	7.187	686	697	716	rVB2	169678	384564	12.33%	0.840%
7	7.499	744	750	762	rBV	104262	179321	5.75%	0.392%
8	7.611	762	769	772	rBV	315523	518681	16.63%	1.134%
9	7.652	772	776	786	rVB	400460	698206	22.39%	1.526%
10	7.958	820	828	840	rBV	272673	497438	15.95%	1.087%
11	8.346	887	894	901	rBV	34074	75496	2.42%	0.165%
12	8.422	901	907	926	rVB	203030	391542	12.55%	0.856%
13	8.663	940	948	955	rVB	251096	429692	13.78%	0.939%
14	8.781	962	968	972	rBV	37682	71039	2.28%	0.155%
15	8.828	972	976	989	rVB	62250	129352	4.15%	0.283%
16	9.522	1082	1094	1108	rBV2	73216	216010	6.93%	0.472%
17	9.822	1136	1145	1161	rBV	93308	211297	6.77%	0.462%
18	10.069	1177	1187	1210	rBV6	87473	326464	10.47%	0.713%
19	10.381	1232	1240	1244	rBV	381844	743254	23.83%	1.624%
20	10.434	1244	1249	1264	rVV	513109	984008	31.55%	2.151%
21	10.581	1268	1274	1286	rBV2	20588	56521	1.81%	0.124%
22	10.799	1308	1311	1317	rVB8	2149	3673	0.12%	0.008%
23	11.199	1371	1379	1382	rBV8	2674	6462	0.21%	0.014%
24	11.704	1458	1465	1485	rBV	223700	579253	18.57%	1.266%
25	12.534	1600	1606	1609	rBV7	1746	3238	0.10%	0.007%
26	12.693	1626	1633	1641	rBV	235374	396976	12.73%	0.868%
27	12.775	1641	1647	1672	rVB	360573	851472	27.30%	1.861%
28	13.134	1702	1708	1719	rBV	275722	480572	15.41%	1.050%
29	13.328	1737	1741	1744	rVB6	2758	3571	0.11%	0.008%
30	13.516	1765	1773	1778	rBV6	2407	3995	0.13%	0.009%
31	13.704	1798	1805	1823	rBV	109010	201770	6.47%	0.441%
32	13.881	1829	1835	1845	rBV	650204	1078655	34.58%	2.357%
33	13.998	1845	1855	1869	rVB2	951950	1568739	50.30%	3.428%
34	14.281	1894	1903	1908	rBV	746142	1020203	32.71%	2.230%
35	14.345	1908	1914	1923	rVV	759567	1122953	36.00%	2.454%
36	14.422	1924	1927	1931	rVB	12555	14970	0.48%	0.033%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Title : SVOA CALIBRATION

37	14.510	1933	1942	1948	rBV	590095	1381005	44.28%	3.018%
38	14.581	1948	1954	1969	rVV2	155884	506681	16.24%	1.107%
39	14.704	1970	1975	1991	rVB	362163	677849	21.73%	1.481%
40	14.945	2012	2016	2024	rVB7	8861	15433	0.49%	0.034%
41	15.145	2044	2050	2060	rVV	1089746	1568319	50.28%	3.427%
42	15.304	2071	2077	2082	rBV	215855	303461	9.73%	0.663%
43	15.363	2082	2087	2101	rVB	1178173	1738039	55.72%	3.798%
44	15.481	2103	2107	2118	rVV	34220	58211	1.87%	0.127%
45	15.622	2127	2131	2135	rVB6	5858	7992	0.26%	0.017%
46	15.881	2169	2175	2176	rBV5	3051	4872	0.16%	0.011%
47	15.922	2179	2182	2187	rVV2	6964	11027	0.35%	0.024%
48	15.969	2187	2190	2195	rVB7	2902	4501	0.14%	0.010%
49	16.081	2202	2209	2216	rBV7	6736	16844	0.54%	0.037%
50	16.251	2235	2238	2245	rVB5	8557	13746	0.44%	0.030%
51	16.387	2253	2261	2270	rBV	1350340	2050301	65.74%	4.481%
52	16.528	2283	2285	2293	rVB9	2829	4493	0.14%	0.010%
53	16.757	2318	2324	2338	rBV	679675	1198459	38.42%	2.619%
54	17.116	2378	2385	2389	rBV	843479	1277920	40.97%	2.793%
55	17.157	2389	2392	2398	rVV	452664	669148	21.45%	1.462%
56	17.216	2398	2402	2405	rVV	216568	344974	11.06%	0.754%
57	17.257	2405	2409	2423	rVB	544708	930294	29.83%	2.033%
58	17.763	2491	2495	2500	rBV5	15094	24126	0.77%	0.053%
59	17.839	2504	2508	2514	rVB	31370	44350	1.42%	0.097%
60	18.010	2535	2537	2542	rVB6	2988	4310	0.14%	0.009%
61	18.075	2542	2548	2554	rBV2	40312	75900	2.43%	0.166%
62	18.145	2554	2560	2569	rVV	1230207	1721379	55.19%	3.762%
63	18.369	2594	2598	2603	rBV7	5400	11195	0.36%	0.024%
64	18.475	2613	2616	2621	rBV7	8770	16536	0.53%	0.036%
65	18.628	2636	2642	2652	rVB2	27621	66131	2.12%	0.145%
66	18.980	2698	2702	2708	rBV9	12003	23651	0.76%	0.052%
67	19.069	2713	2717	2721	rVB6	25888	34054	1.09%	0.074%
68	19.222	2739	2743	2745	rBV3	21336	24942	0.80%	0.055%
69	19.275	2745	2752	2758	rBV	1603484	2310218	74.07%	5.049%
70	19.416	2773	2776	2784	rVB7	26576	40043	1.28%	0.088%
71	19.533	2792	2796	2799	rBV3	11648	16732	0.54%	0.037%
72	19.628	2806	2812	2813	rBV	292333	398325	12.77%	0.871%
73	19.657	2813	2817	2832	rVB	737678	1139138	36.52%	2.490%
74	20.375	2936	2939	2944	rBV2	29999	38482	1.23%	0.084%
75	21.463	3121	3124	3129	rVB3	51386	65146	2.09%	0.142%
76	21.533	3131	3136	3141	rVV	1872683	2585068	82.88%	5.650%

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

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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Title : SVOA CALIBRATION

77	21.604	3141	3148	3153	rVV2	1293646	3119033	100.00%	6.817%
78	21.657	3153	3157	3167	rVB	1480920	2422102	77.66%	5.293%
79	23.974	3542	3551	3562	rVB	934098	2381240	76.35%	5.204%
80	24.727	3674	3679	3691	rVB	136131	375532	12.04%	0.821%
81	24.963	3711	3719	3731	rVB	514449	1439437	46.15%	3.146%
82	28.833	4368	4377	4378	rBV	308814	639228	20.49%	1.397%

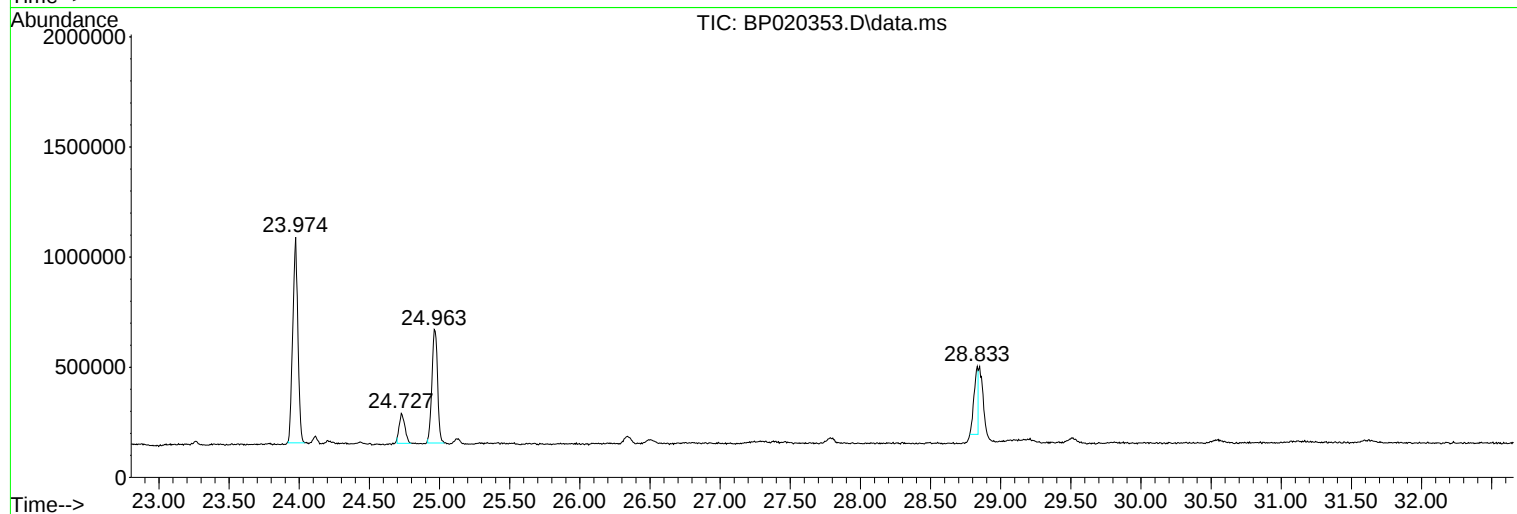
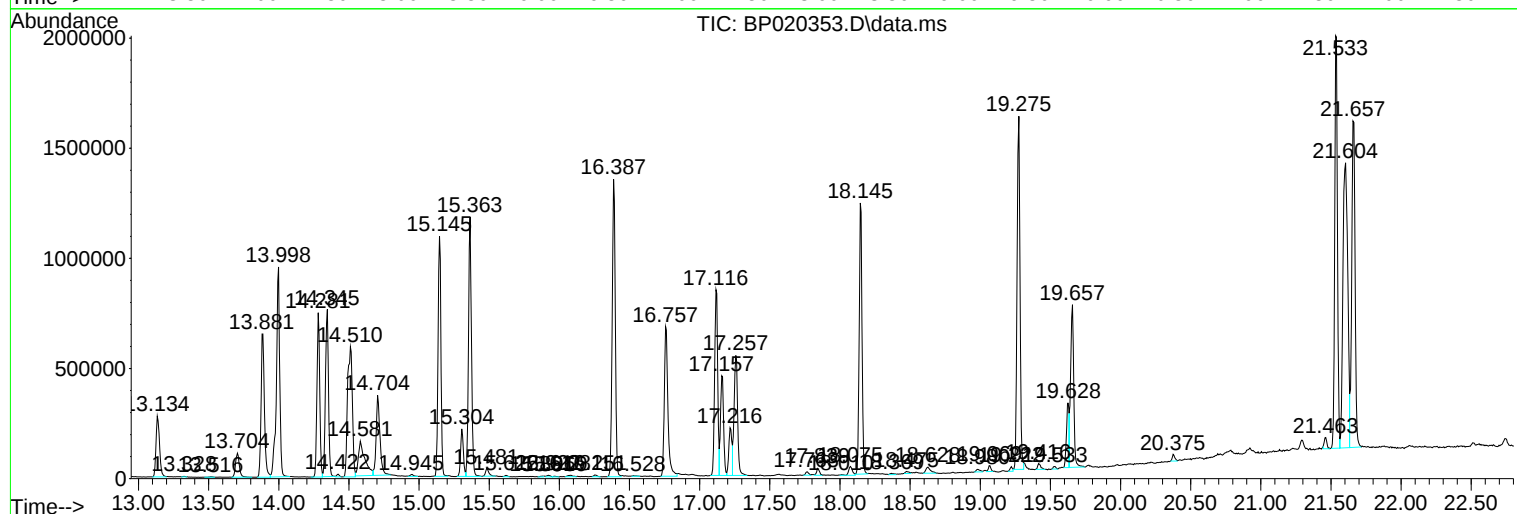
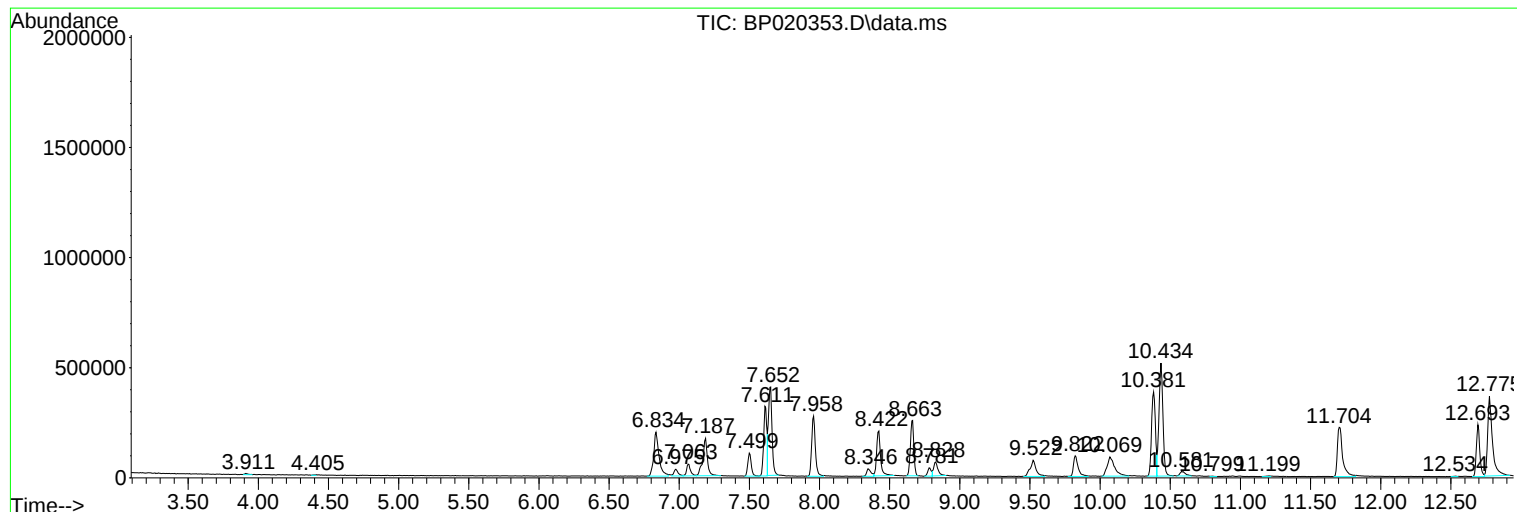
Sum of corrected areas: 45757049

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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



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Sample : P2372-23DL 5X
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ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Y3DL

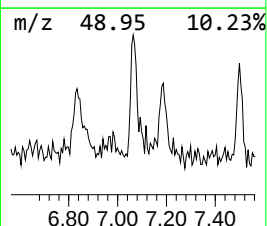
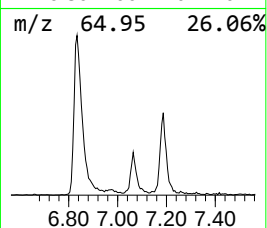
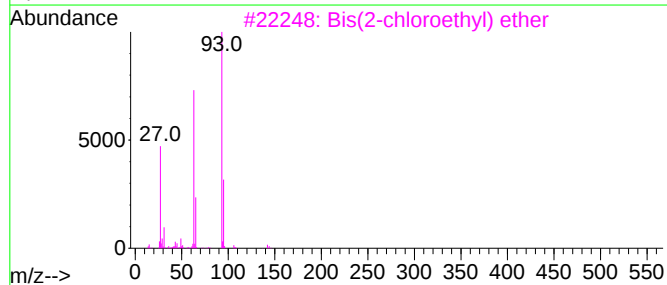
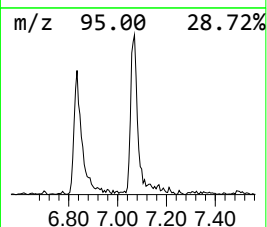
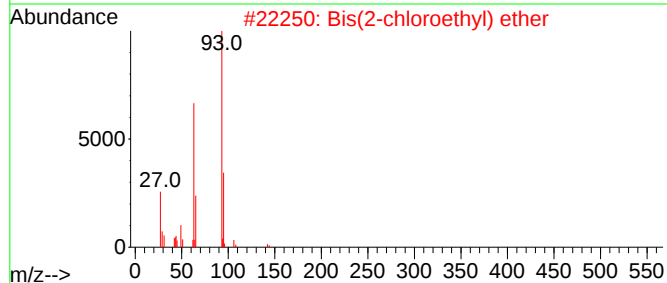
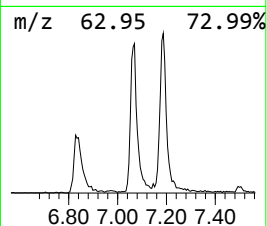
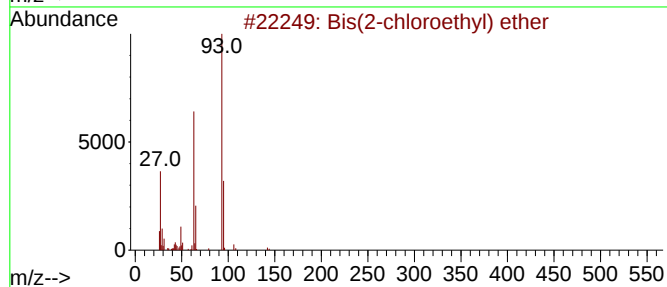
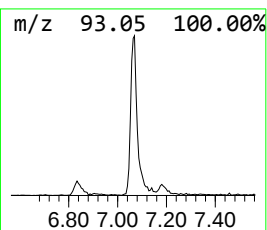
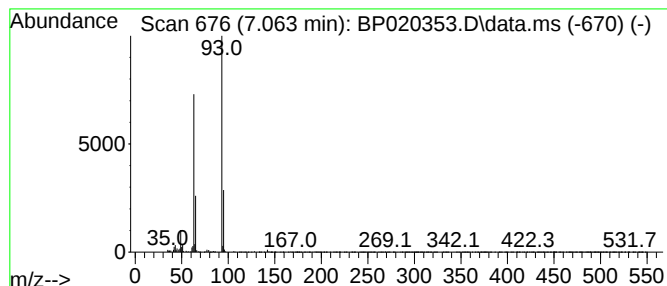
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Bis(2-chloroethyl) ether Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.063	4.09 ng/ul	106097	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	97
2			Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	90
3			Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	86
4			Methane, bis(2-chloroethoxy)-	172	C5H10Cl2O2	000111-91-1	83
5			Methane, bis(2-chloroethoxy)-	172	C5H10Cl2O2	000111-91-1	78



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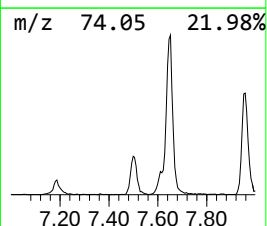
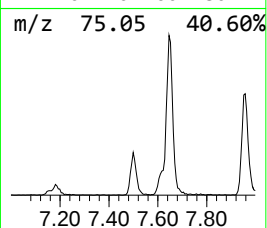
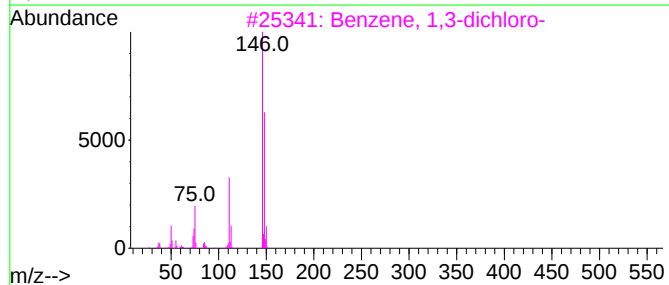
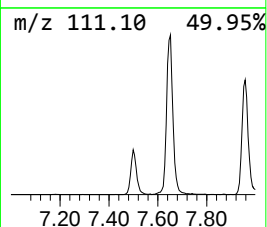
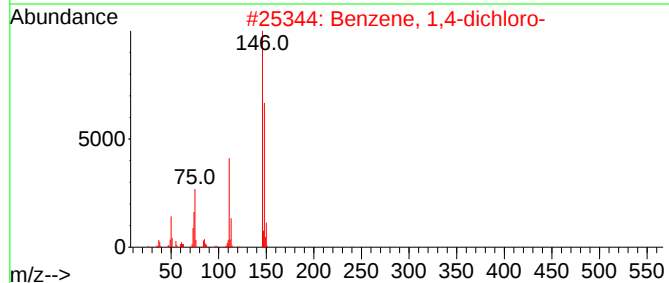
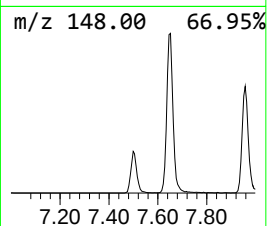
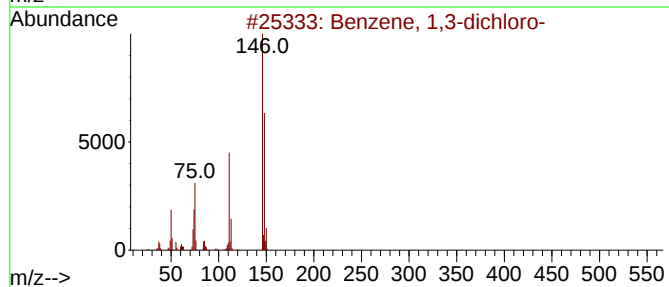
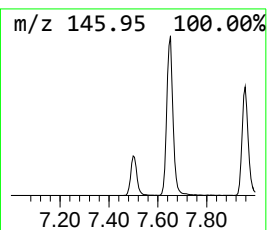
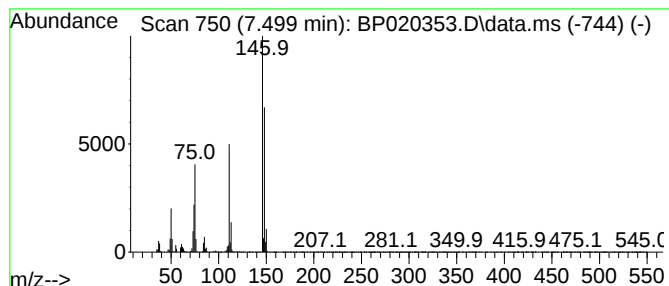
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dichloro- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.499	6.91 ng/ul	179321	1,4-Dichlorobenzene-d4	7.616

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	96
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96



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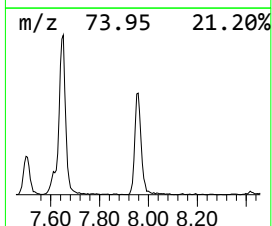
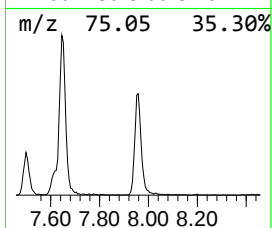
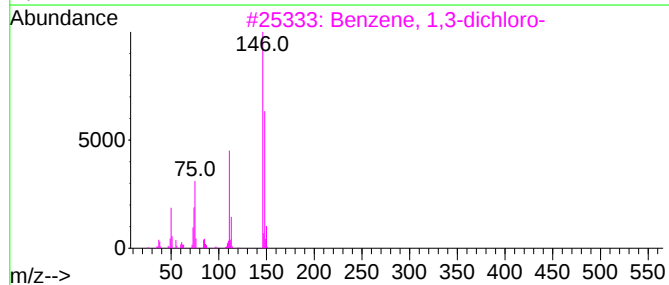
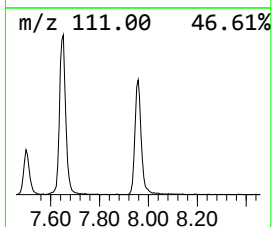
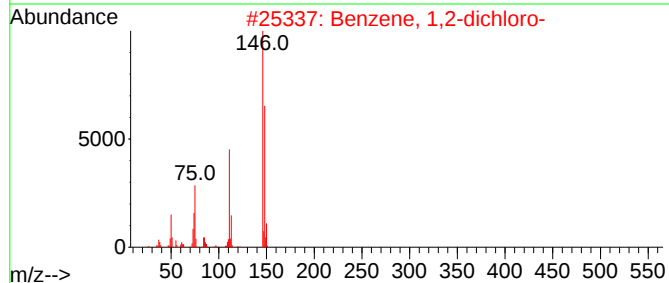
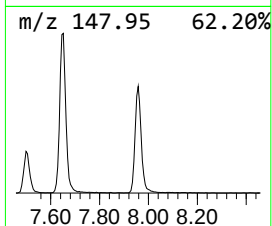
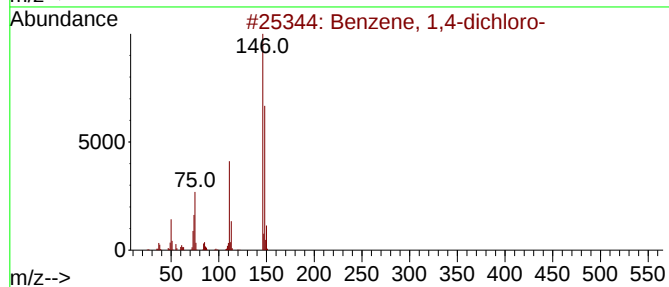
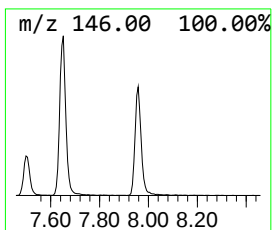
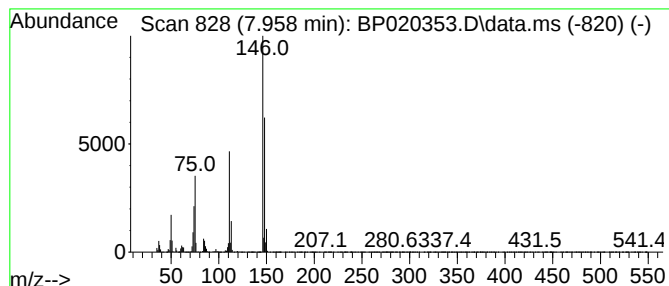
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.958	19.18 ng/ul	497438	1,4-Dichlorobenzene-d4	7.616

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



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A43Y3DL

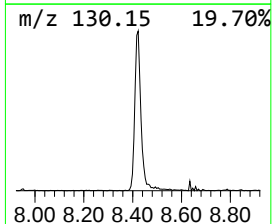
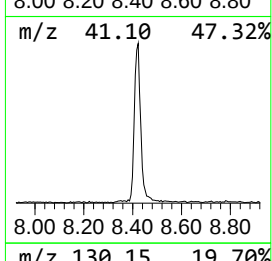
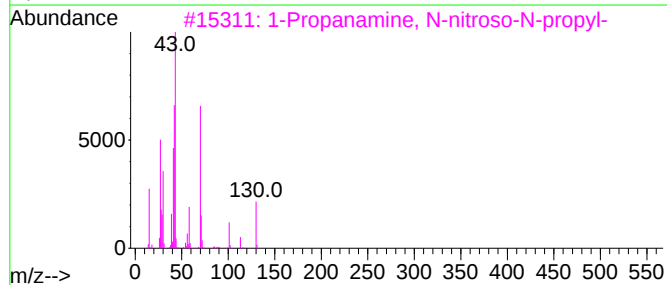
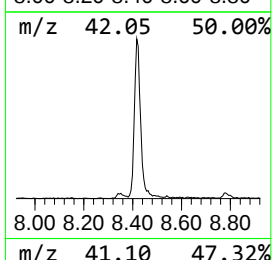
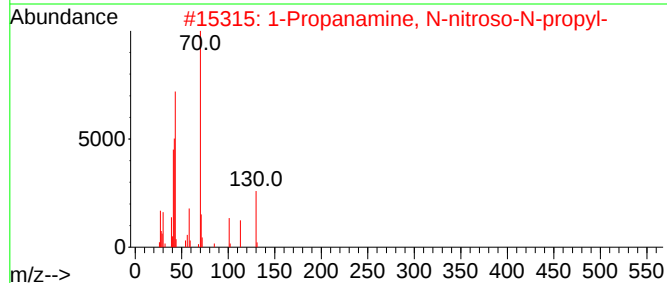
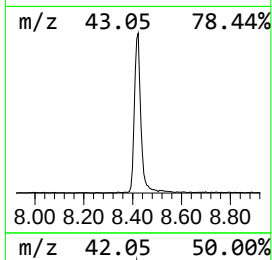
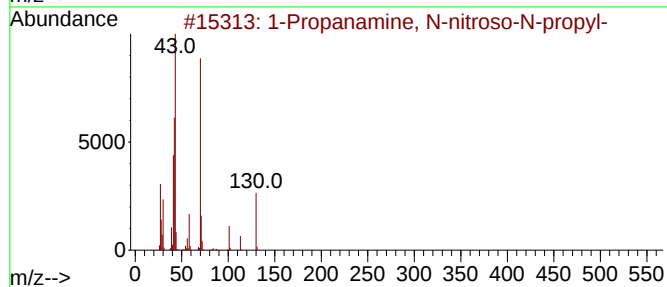
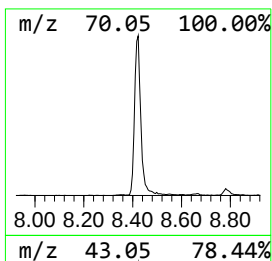
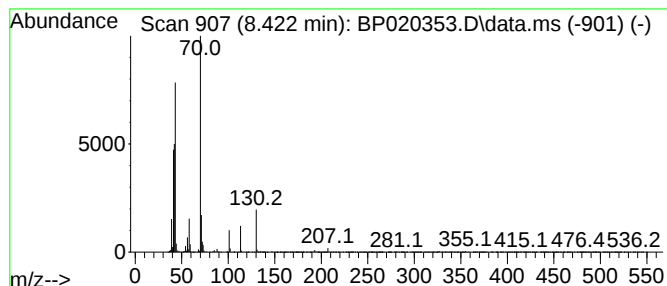
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Propanamine, N-nitroso-N-... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.422	15.10 ng/ul	391542	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	91
2			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	83
3			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	68
4			Oxazolidine, 2,2-dimethyl-	101	C5H11NO	020515-62-2	45
5			1-Propanamine, N-nitroso-N-propyl-	130	C6H14N2O	000621-64-7	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Y3DL

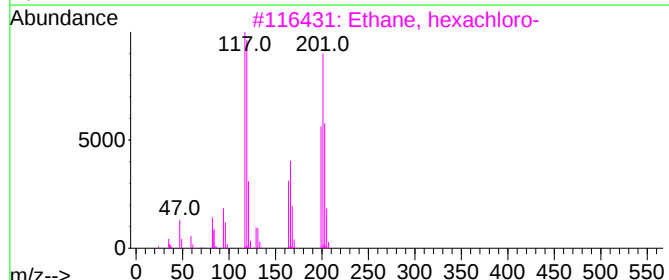
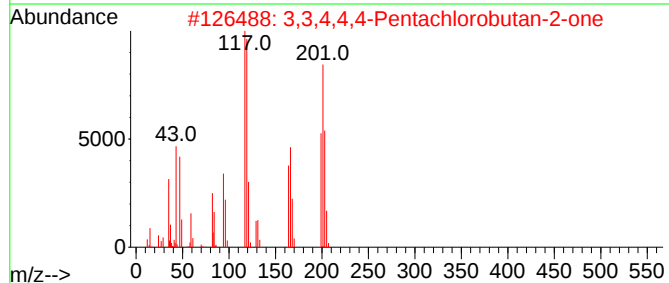
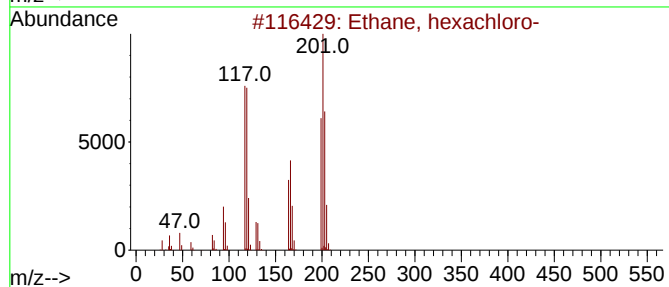
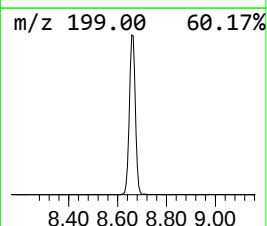
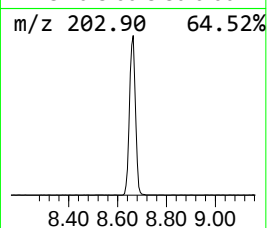
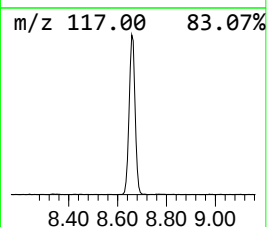
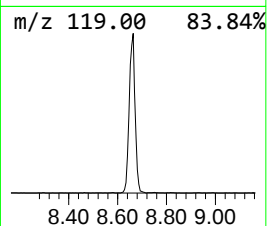
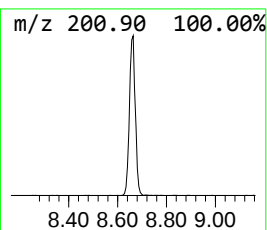
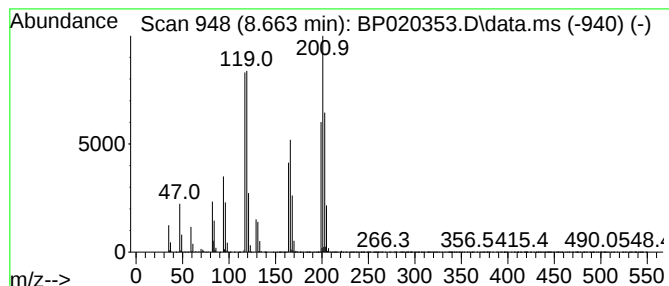
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Ethane, hexachloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.663	16.57 ng/ul	429692	1,4-Dichlorobenzene-d4	7.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, hexachloro-	234	C2Cl6	000067-72-1	91
2			3,3,4,4,4-Pentachlorobutan-2-one	242	C4H3Cl5O	004020-56-8	91
3			Ethane, hexachloro-	234	C2Cl6	000067-72-1	90
4			Ethane, hexachloro-	234	C2Cl6	000067-72-1	87
5			Ethane, hexachloro-	234	C2Cl6	000067-72-1	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Y3DL

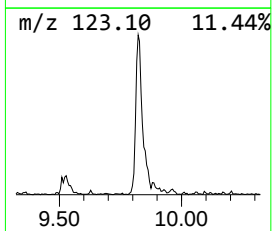
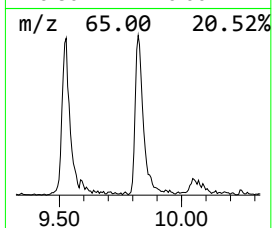
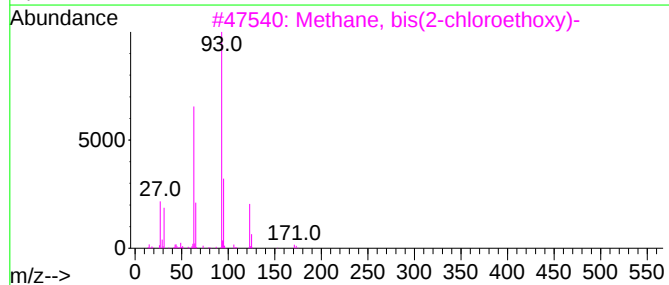
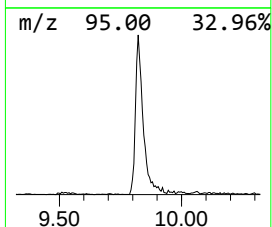
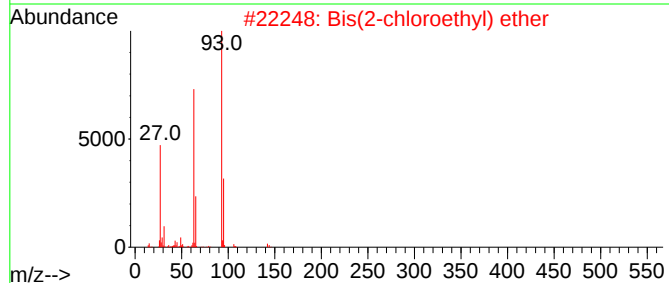
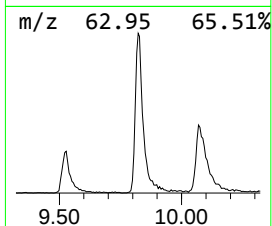
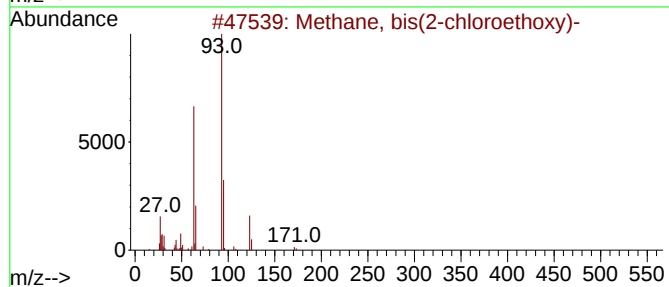
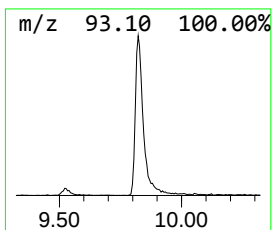
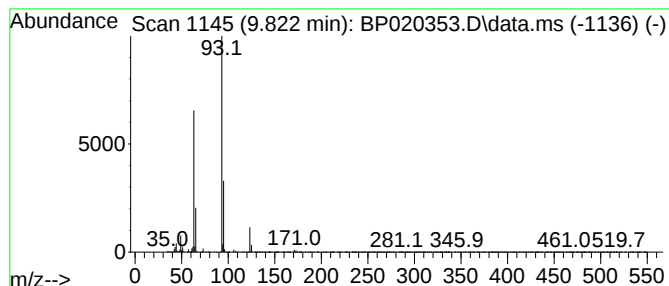
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Methane, bis(2-chloroethoxy)- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.822	5.69 ng/ul	211297	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methane, bis(2-chloroethoxy)-	172	C5H10Cl2O2	000111-91-1	90
2			Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	90
3			Methane, bis(2-chloroethoxy)-	172	C5H10Cl2O2	000111-91-1	74
4			Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	64
5			Bis(2-chloroethyl) ether	142	C4H8Cl2O	000111-44-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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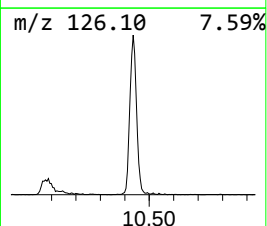
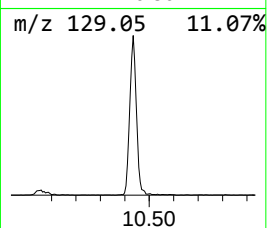
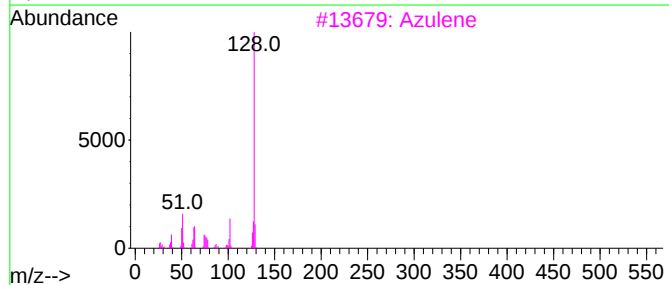
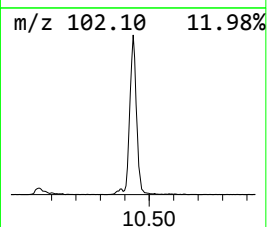
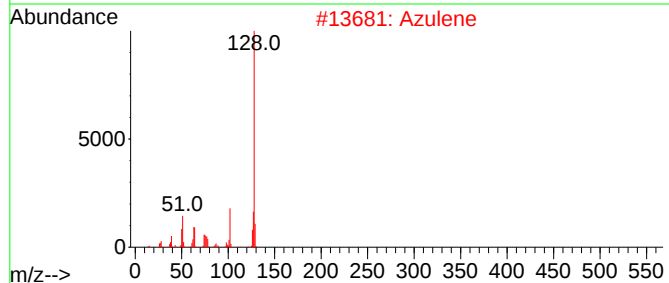
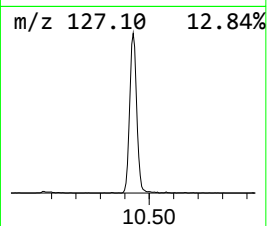
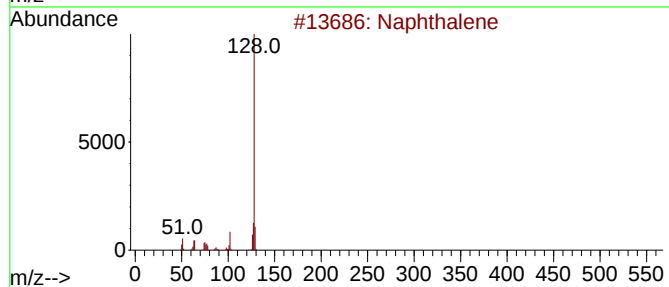
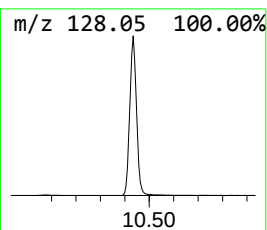
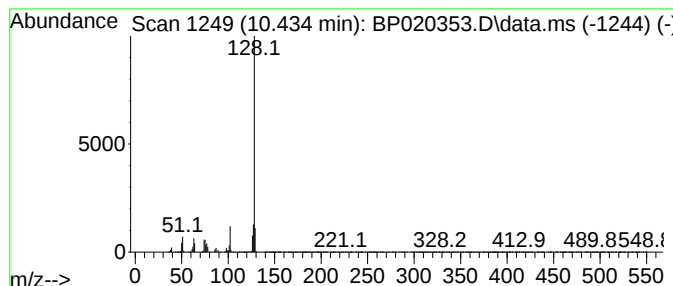
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Naphthalene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.434	26.48 ng/ul	984008	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	91
3			Azulene	128	C10H8	000275-51-4	91
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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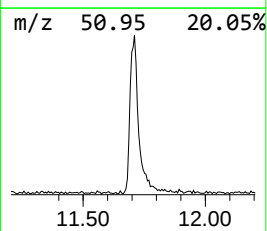
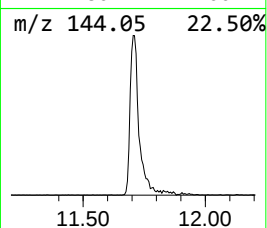
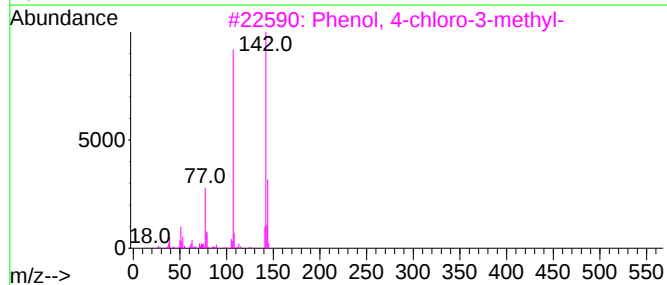
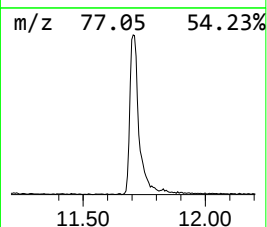
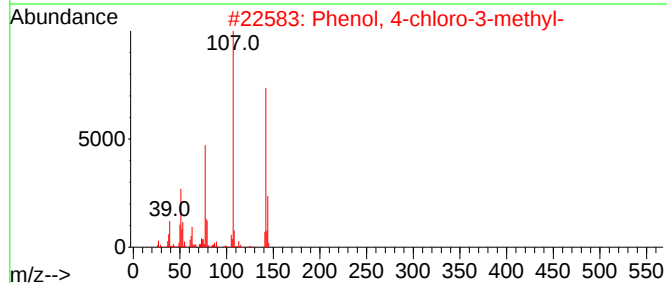
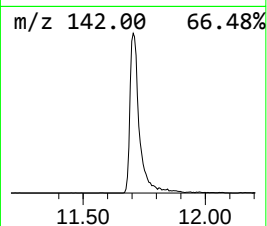
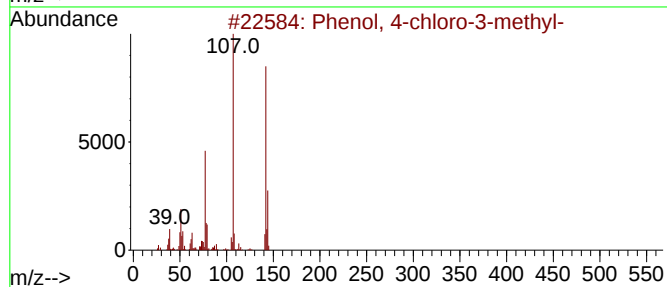
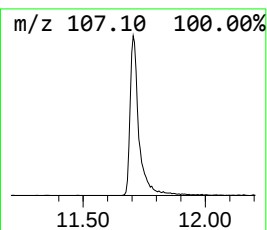
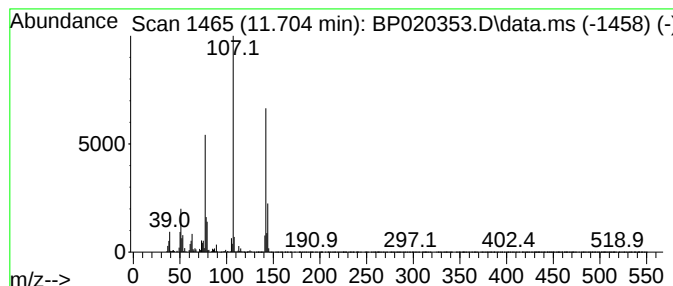
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Phenol, 4-chloro-3-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.704	15.59 ng/ul	579253	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	96
2			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	96
3			Phenol, 4-chloro-3-methyl-	142	C7H7ClO	000059-50-7	95
4			Phenol, 4-chloro-2-methyl-	142	C7H7ClO	001570-64-5	94
5			Phenol, 4-chloro-2-methyl-	142	C7H7ClO	001570-64-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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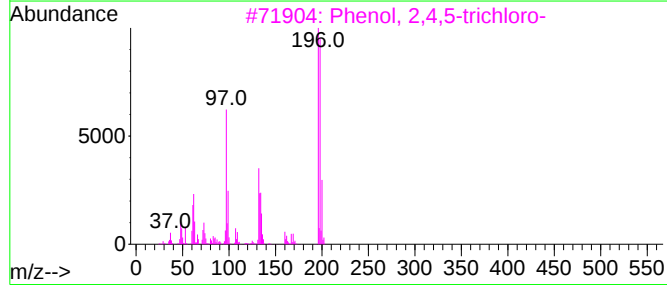
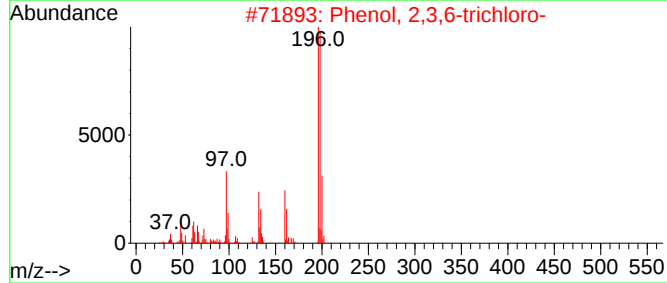
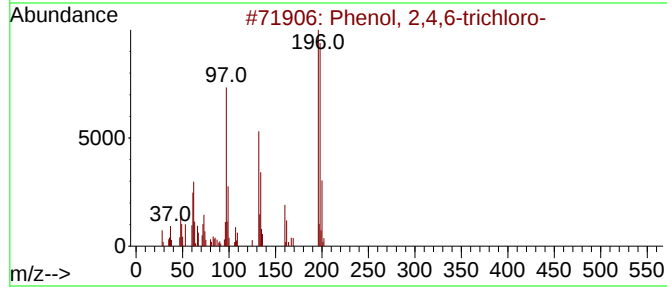
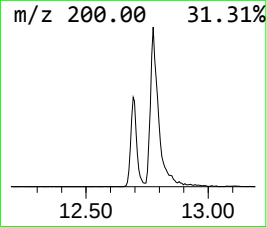
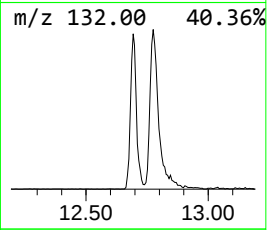
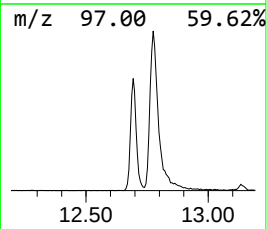
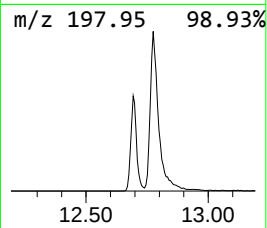
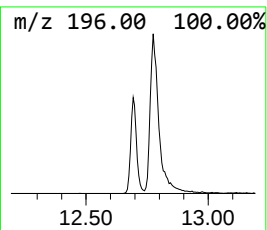
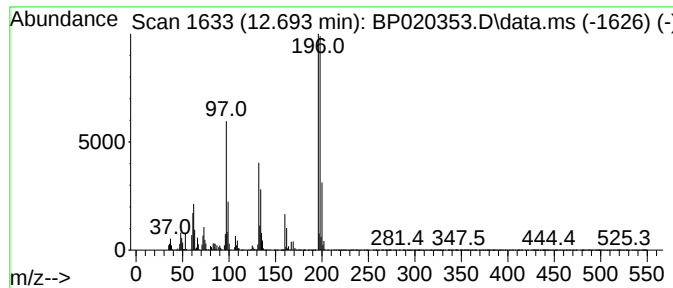
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenol, 2,4,6-trichloro- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.693	7.78 ng/ul	396976	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	99
2			Phenol, 2,3,6-trichloro-	196	C6H3Cl3O	000933-75-5	95
3			Phenol, 2,4,5-trichloro-	196	C6H3Cl3O	000095-95-4	94
4			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	94
5			Phenol, 2,4,6-trichloro-	196	C6H3Cl3O	000088-06-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

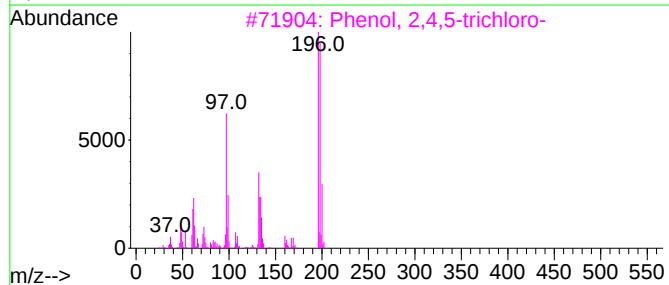
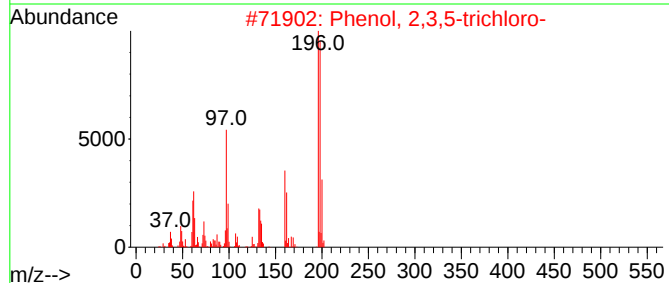
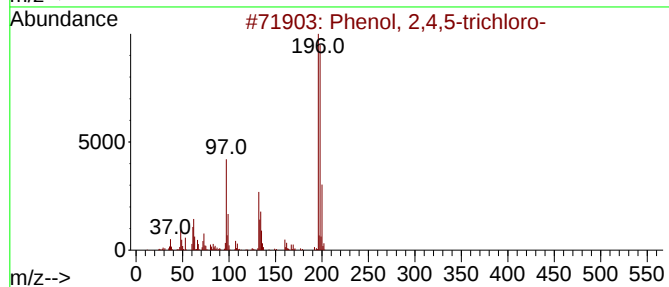
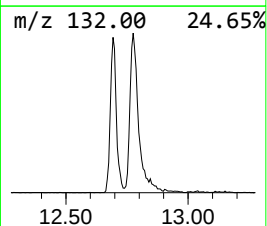
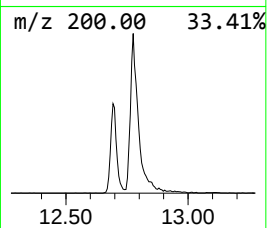
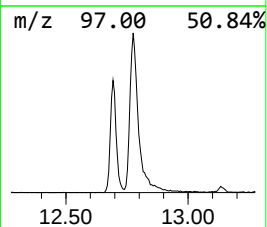
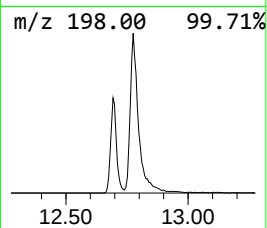
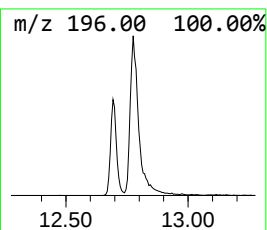
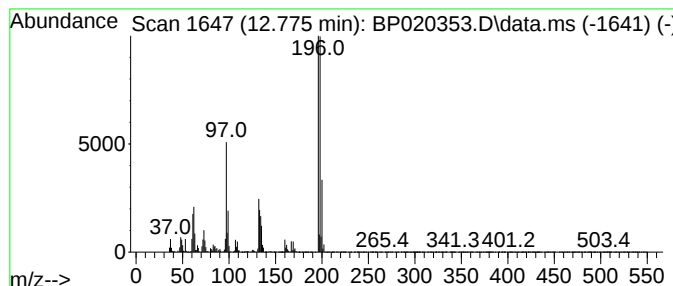
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ClientSampleId :
A43Y3DL

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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Phenol, 2,4,5-trichloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.775	16.69 ng/ul	851472	Acenaphthene-d10			14.281
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Phenol, 2,4,5-trichloro-		196	C6H3Cl3O	000095-95-4	94
2	Phenol, 2,3,5-trichloro-		196	C6H3Cl3O	000933-78-8	94
3	Phenol, 2,4,5-trichloro-		196	C6H3Cl3O	000095-95-4	94
4	Phenol, 2,4,5-trichloro-		196	C6H3Cl3O	000095-95-4	94
5	Phenol, 2,4,6-trichloro-		196	C6H3Cl3O	000088-06-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A43Y3DL

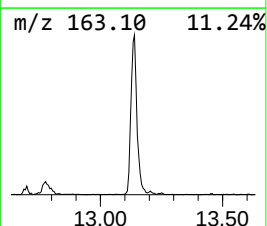
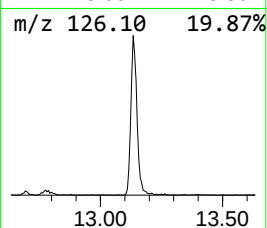
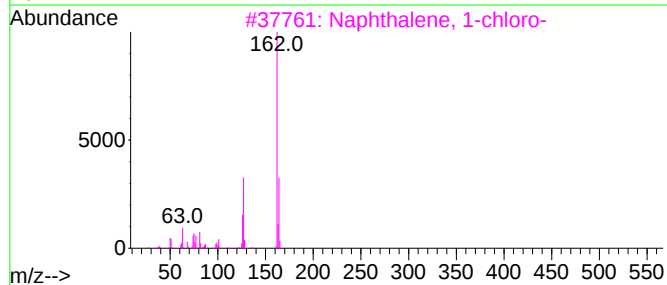
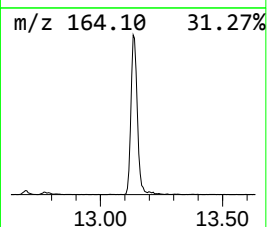
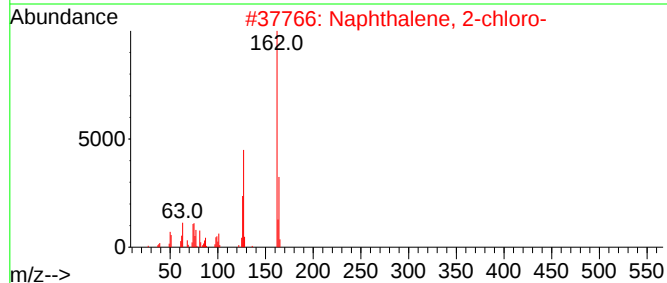
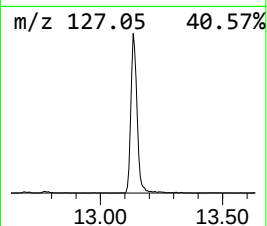
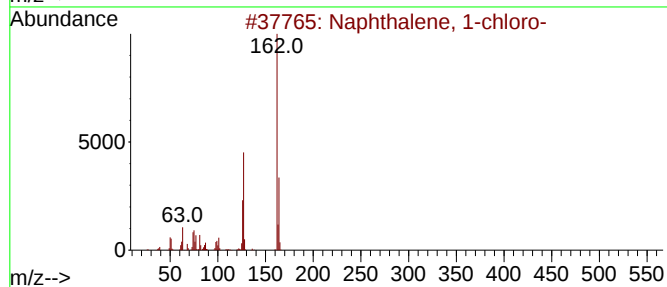
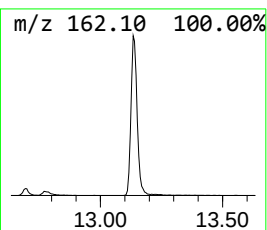
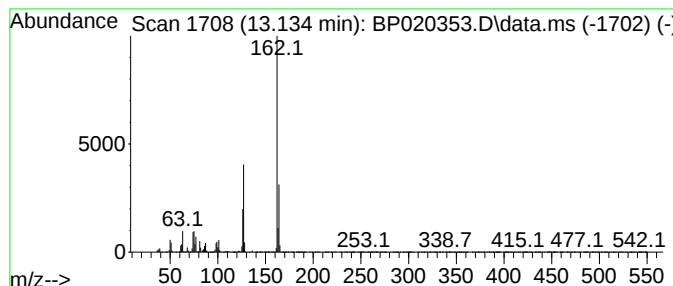
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Naphthalene, 1-chloro- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.134	9.42 ng/ul	480572	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	97
2			Naphthalene, 2-chloro-	162	C10H7Cl	000091-58-7	96
3			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	95
4			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	94
5			Naphthalene, 1-chloro-	162	C10H7Cl	000090-13-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

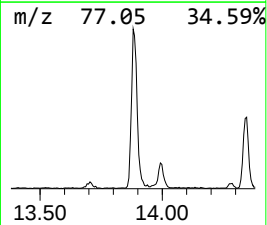
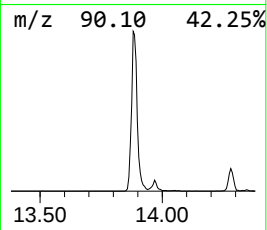
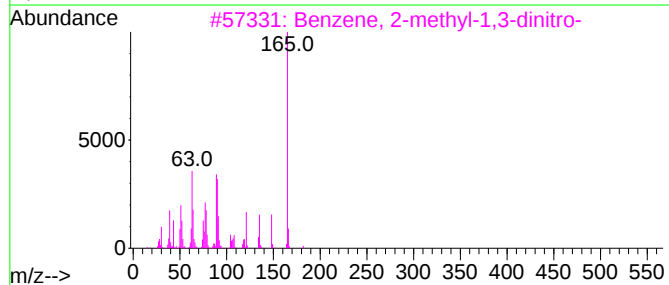
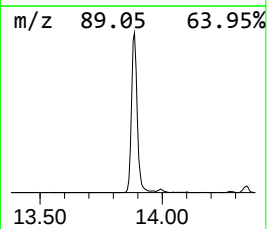
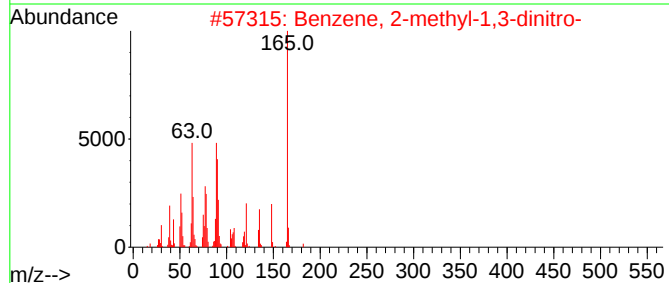
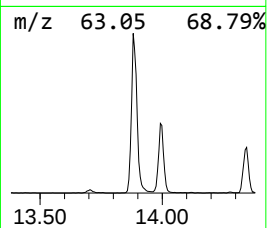
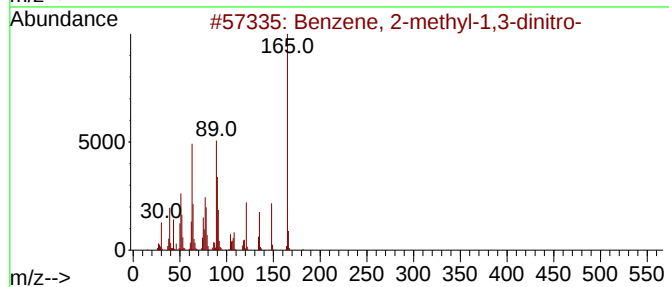
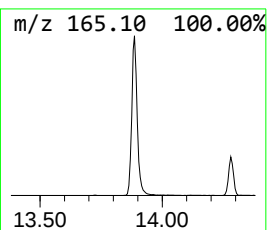
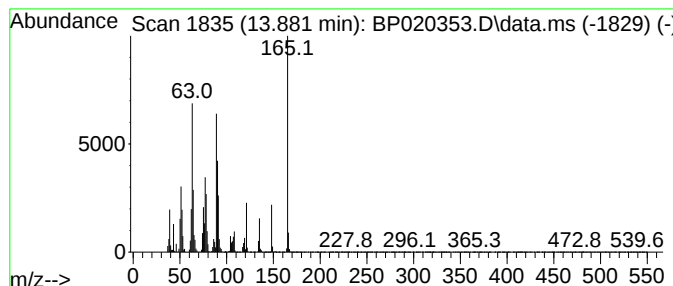
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 2-methyl-1,3-dinitro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.881	21.15 ng/ul	1078660	Acenaphthene-d10			14.281
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 2-methyl-1,3-dinitro-		182	C7H6N2O4	000606-20-2	68
2	Benzene, 2-methyl-1,3-dinitro-		182	C7H6N2O4	000606-20-2	64
3	Benzene, 2-methyl-1,3-dinitro-		182	C7H6N2O4	000606-20-2	59
4	Benzene, 2-methyl-1,3-dinitro-		182	C7H6N2O4	000606-20-2	58
5	1H-Benzotriazole, 4-nitro-		164	C6H4N4O2	006299-39-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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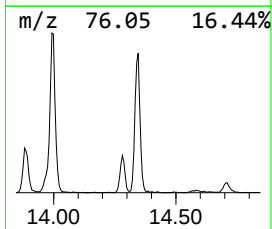
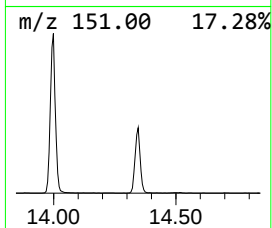
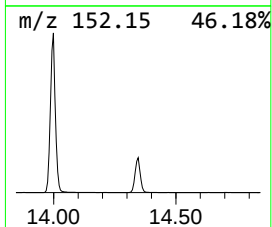
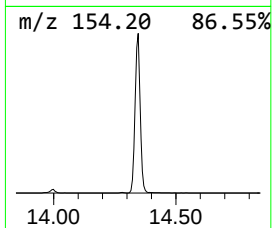
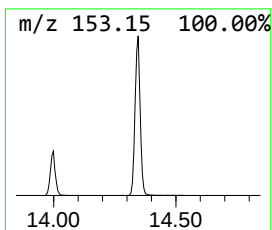
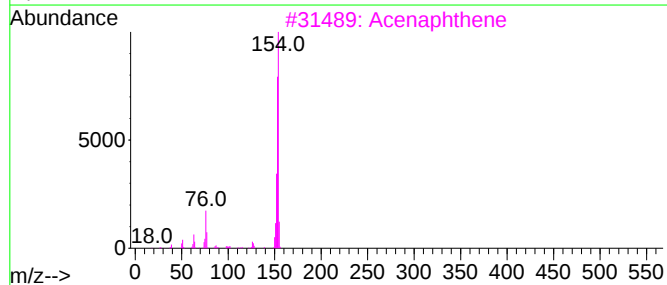
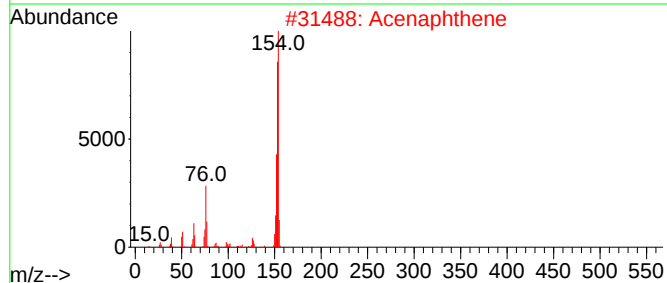
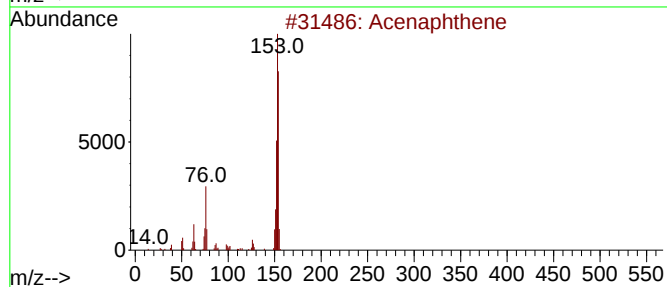
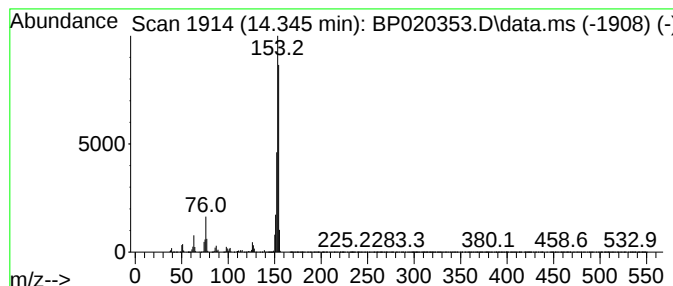
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Acena phthene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.345	22.01 ng/ul	1122950	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthene	154	C12H10	000083-32-9	89
2			Acenaphthene	154	C12H10	000083-32-9	68
3			Acenaphthene	154	C12H10	000083-32-9	64
4			Acenaphthene	154	C12H10	000083-32-9	64
5			Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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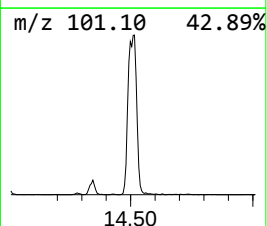
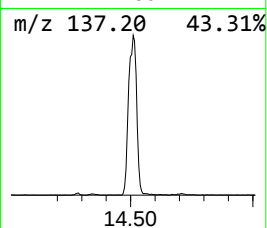
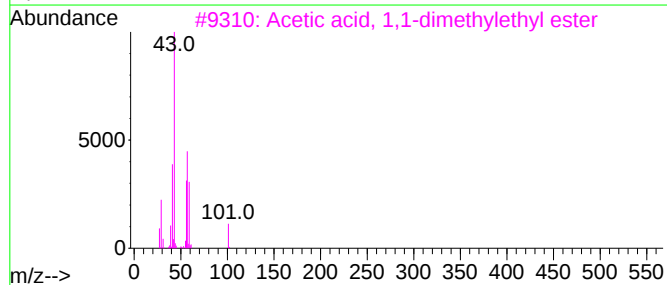
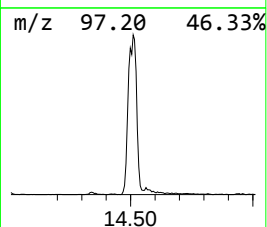
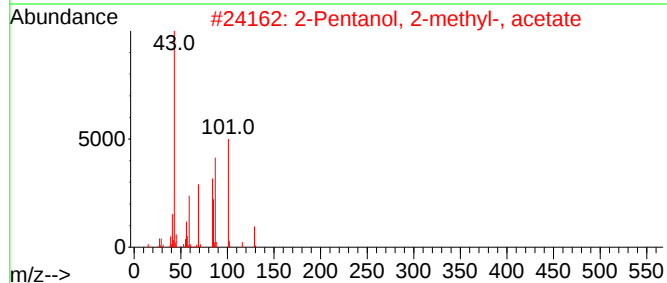
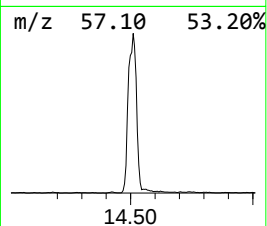
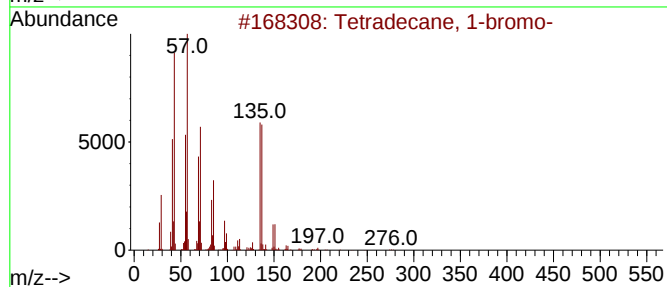
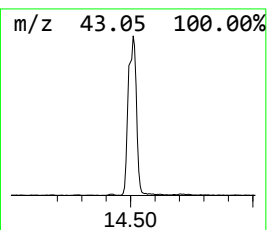
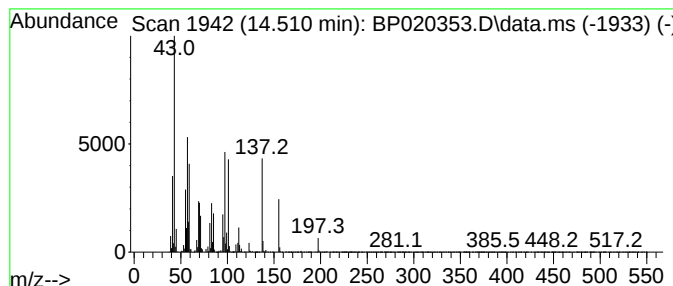
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.510	27.07 ng/ul	1381010	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	14
2			2-Pentanol, 2-methyl-, acetate	144	C8H16O2	034859-98-8	14
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	14
4			trans-5-Isopropyl-6,7-epoxy-8-hy...	228	C13H24O3	058002-07-6	12
5			2-Piperidinone, N-[4-bromo-n-but...	233	C9H16BrNO	195194-80-0	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

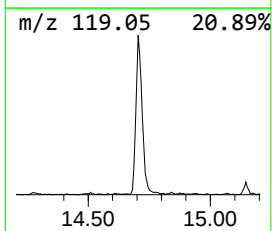
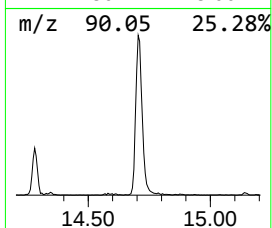
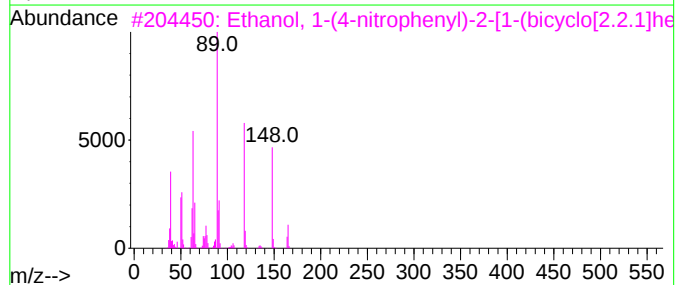
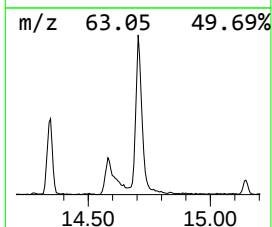
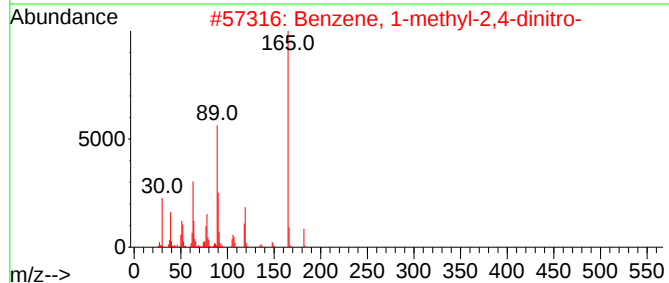
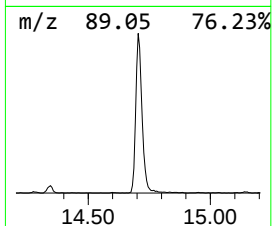
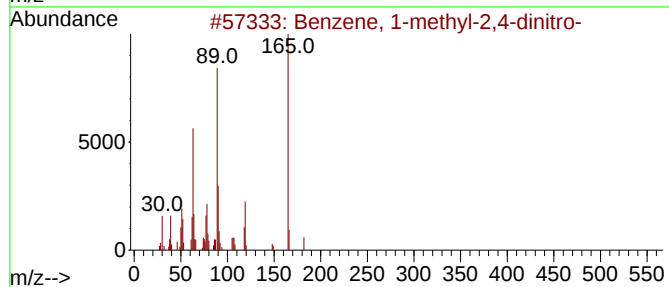
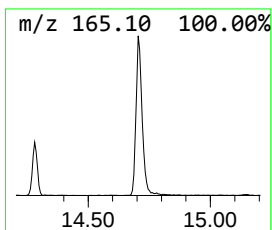
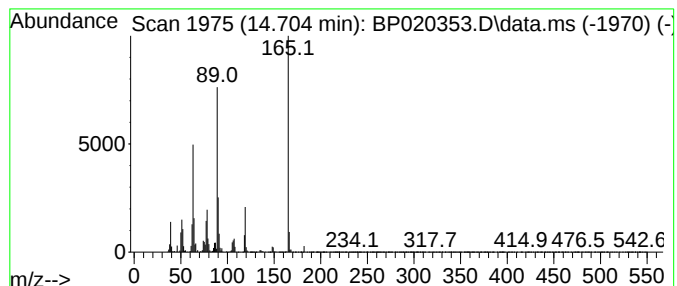
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1-methyl-2,4-dinitro- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.704	13.29 ng/ul	677849	Acenaphthene-d10			14.281
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2,4-dinitro-		182	C7H6N2O4	000121-14-2	91
2	Benzene, 1-methyl-2,4-dinitro-		182	C7H6N2O4	000121-14-2	64
3	Ethanol, 1-(4-nitrophenyl)-2-[1-...		304	C17H24N2O3	1000264-75-7	37
4	Methyl glycolate, TMS derivative		162	C6H14O3Si	1000366-86-9	32
5	6-Nitro-o-tolunitrile		162	C8H6N2O2	001885-76-3	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
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Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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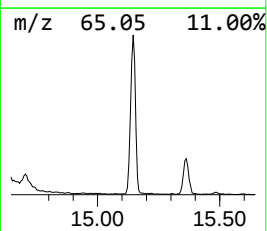
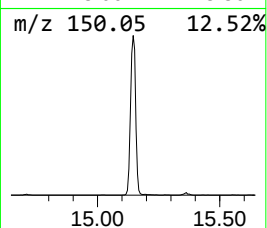
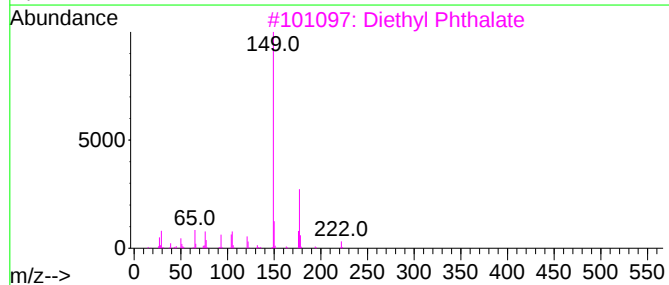
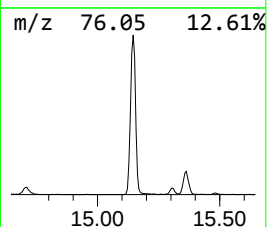
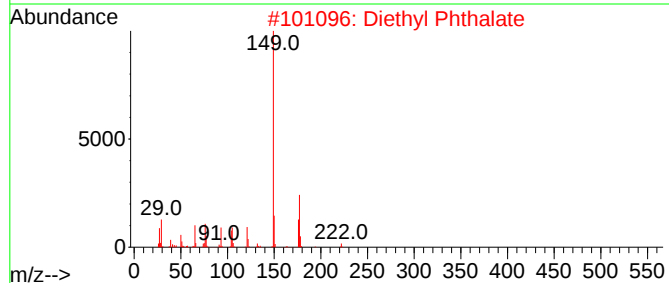
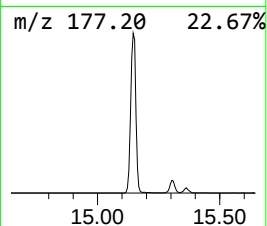
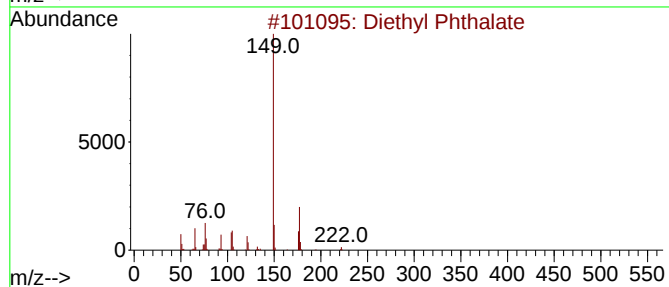
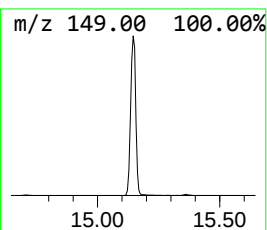
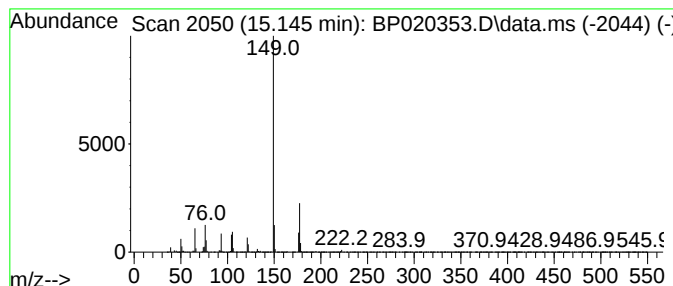
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Diethyl Phthalate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.145	30.75 ng/ul	1568320	Acenaphthene-d10	14.281

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Diethyl Phthalate	222	C12H14O4	000084-66-2	98
2			Diethyl Phthalate	222	C12H14O4	000084-66-2	97
3			Diethyl Phthalate	222	C12H14O4	000084-66-2	93
4			Diethyl Phthalate	222	C12H14O4	000084-66-2	91
5			Phthalic acid, 7-bromoheptyl eth...	370	C17H23BrO4	1000415-51-6	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

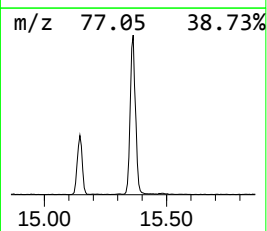
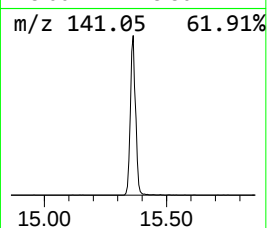
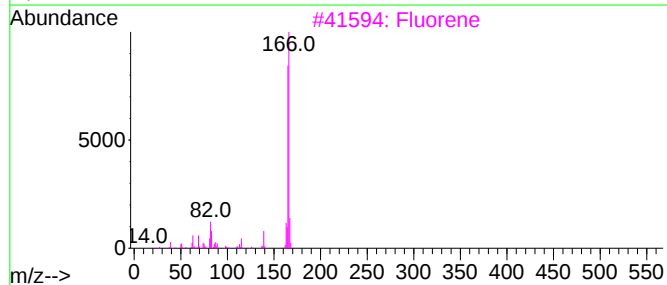
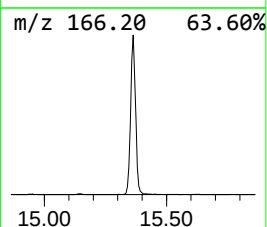
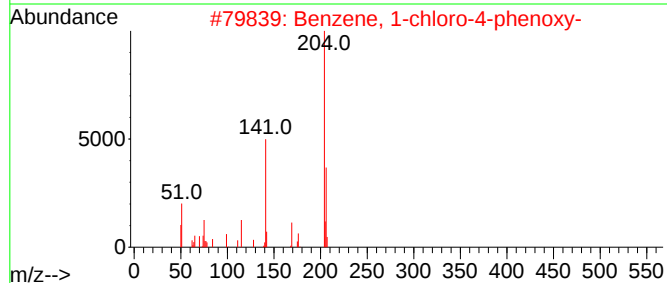
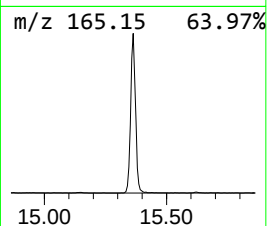
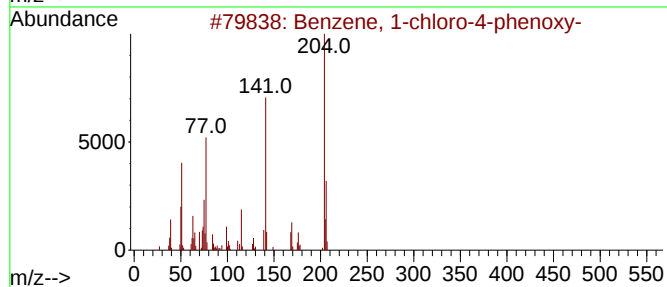
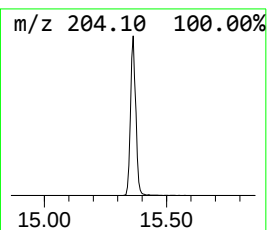
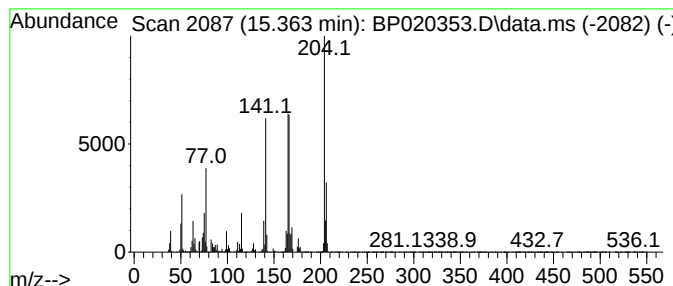
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BNA_P
ClientSampleId :
A43Y3DL

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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzene, 1-chloro-4-phenoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.363	34.07 ng/ul	1738040	Acenaphthene-d10	14.281	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	96
2	Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	96
3	Fluorene	166	C13H10	000086-73-7	62
4	Benzene, 1-chloro-3-phenoxy-	204	C12H9ClO	006452-49-9	60
5	Benzene, 1-chloro-4-phenoxy-	204	C12H9ClO	007005-72-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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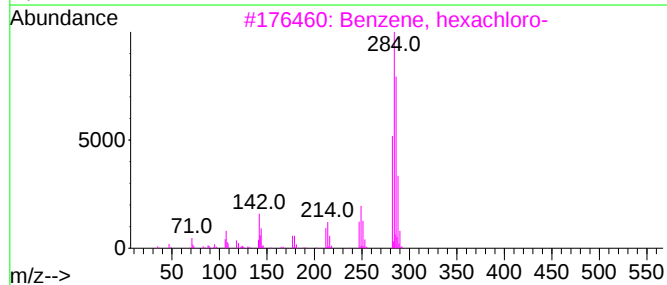
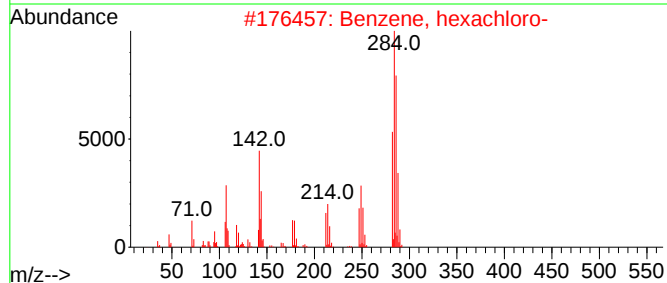
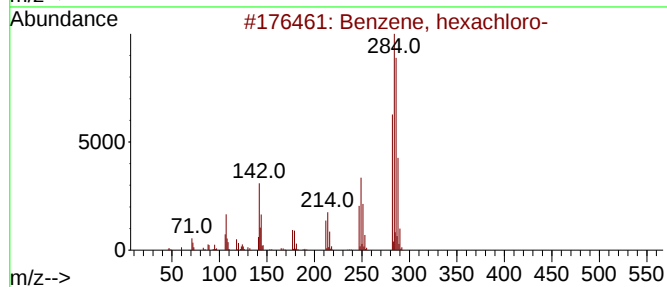
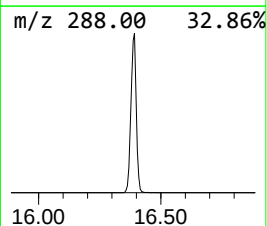
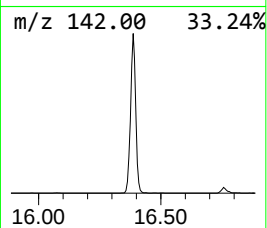
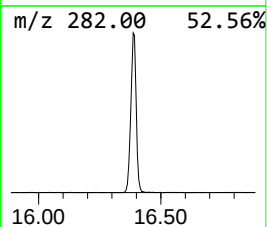
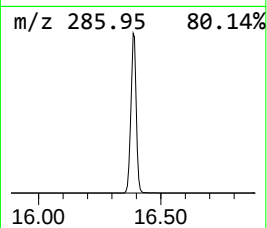
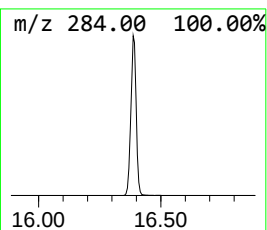
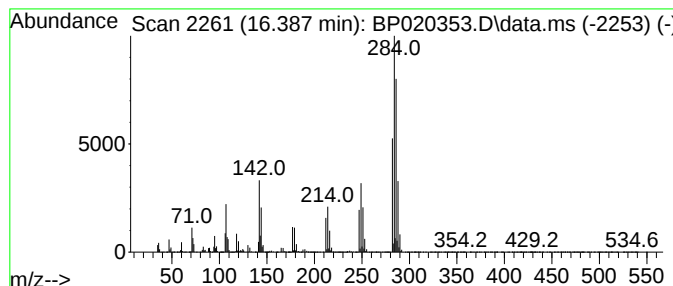
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 Benzene, hexachloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.387	32.09 ng/ul	2050300	Phenanthrene-d10	17.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, hexachloro-	282	C6Cl6	000118-74-1	99
2			Benzene, hexachloro-	282	C6Cl6	000118-74-1	99
3			Benzene, hexachloro-	282	C6Cl6	000118-74-1	96
4			Benzene, hexachloro-	282	C6Cl6	000118-74-1	96
5			Benzene, hexachloro-	282	C6Cl6	000118-74-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

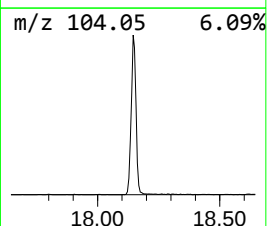
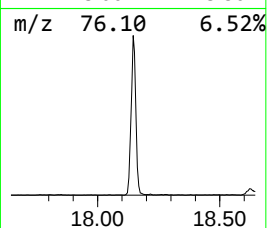
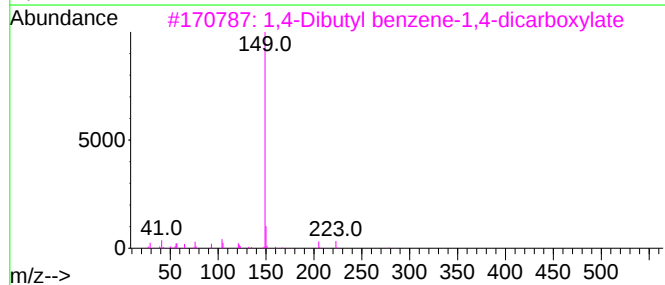
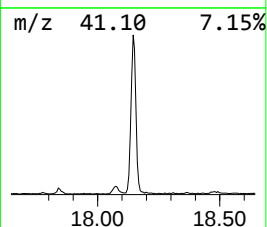
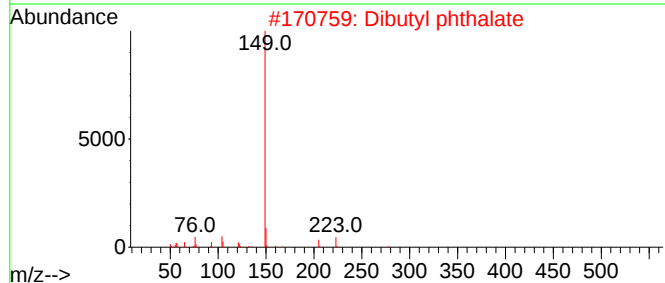
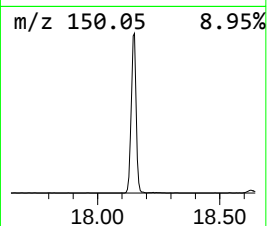
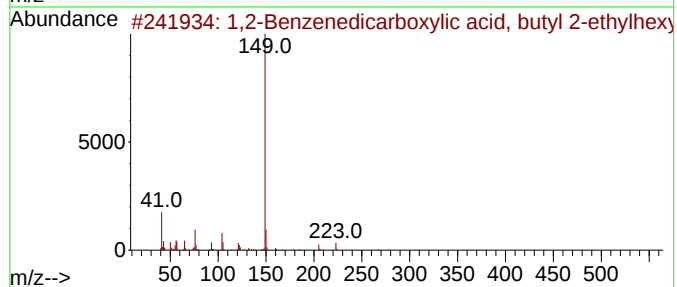
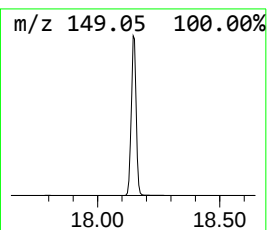
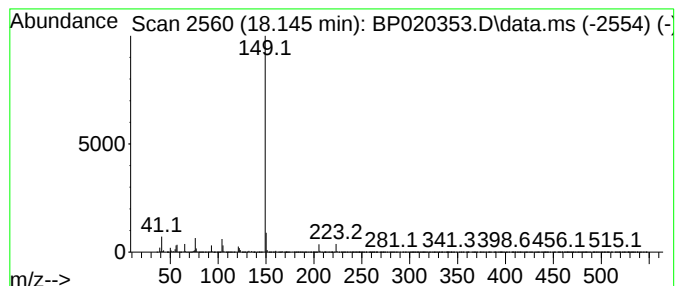
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 19 1,2-Benzenedicarboxylic aci... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.145	26.94 ng/ul	1721380	Phenanthrene-d10	17.116	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	91
2	Dibutyl phthalate	278	C16H22O4	000084-74-2	91
3	1,4-Dibutyl benzene-1,4-dicarbox...	278	C16H22O4	001962-75-0	90
4	Di-sec-butyl phthalate	278	C16H22O4	1000373-65-4	87
5	1,2-Benzenedicarboxylic acid, bu...	304	C18H24O4	000084-64-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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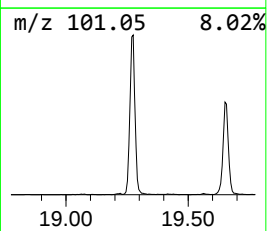
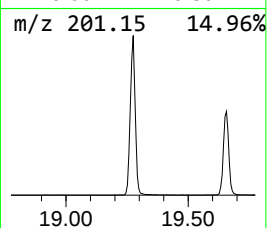
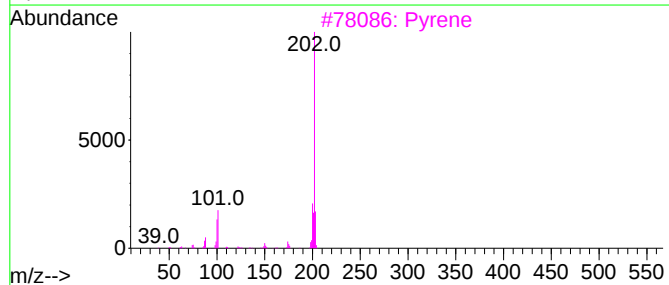
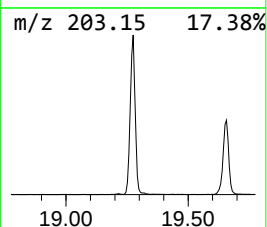
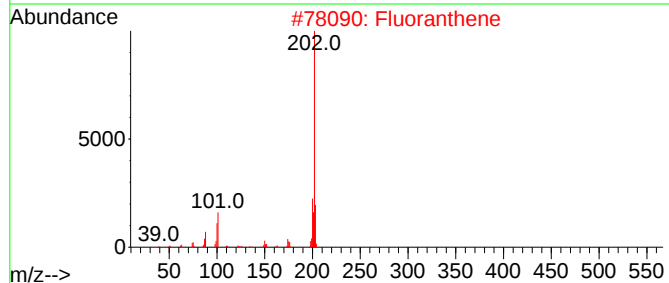
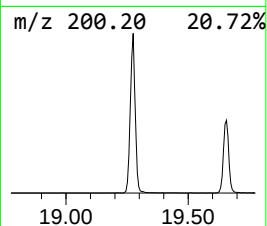
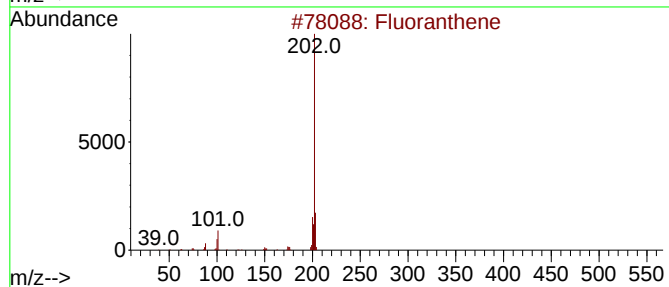
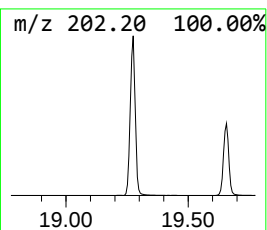
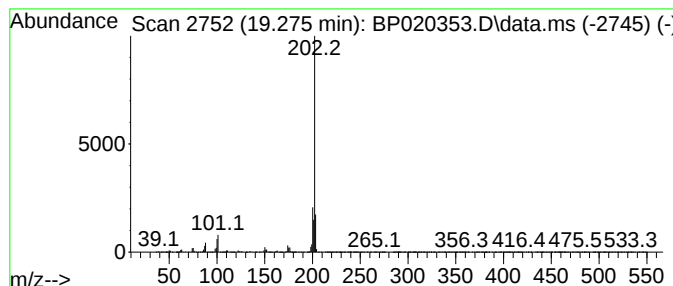
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Fluoranthene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.275	36.16 ng/ul	2310220	Phenanthrene-d10	17.116

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Fluoranthene	202	C16H10	000206-44-0	96
2		Fluoranthene	202	C16H10	000206-44-0	94
3		Pyrene	202	C16H10	000129-00-0	90
4		Pyrene	202	C16H10	000129-00-0	87
5		Pyrene	202	C16H10	000129-00-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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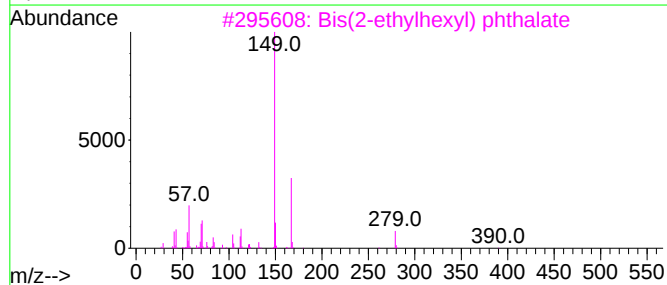
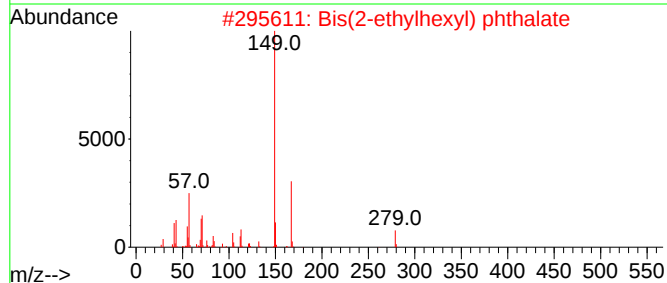
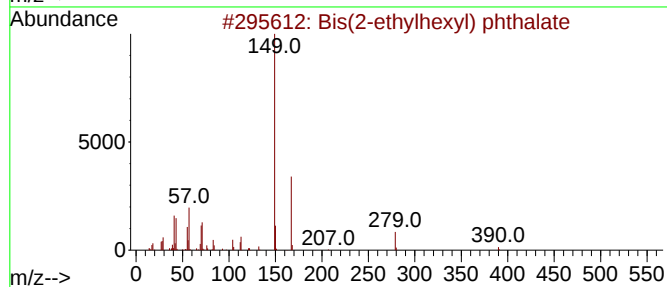
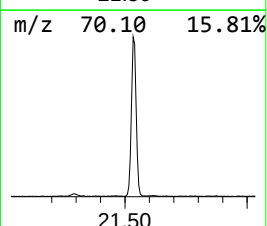
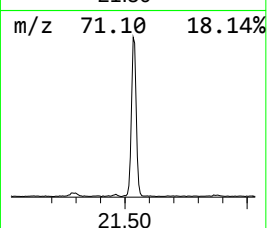
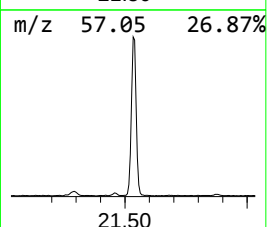
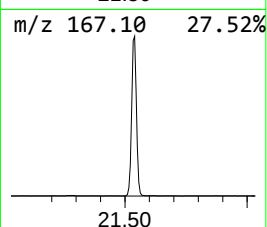
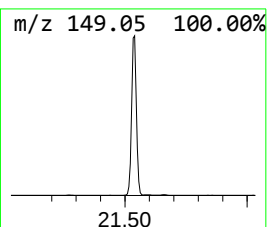
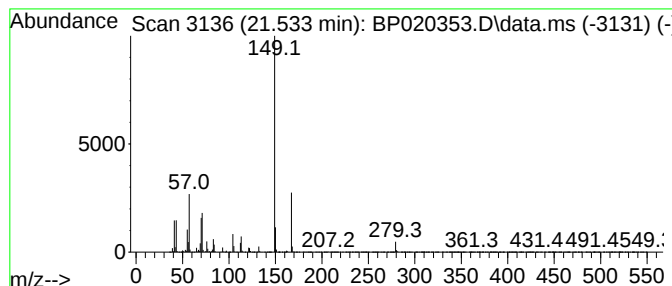
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 Bis(2-ethylhexyl) phthalate Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.533	16.58 ng/ul	2585070	Chrysene-d12	21.616

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	93
2			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	91
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	90
4			Phthalic acid, di(oct-3-yl) ester	390	C24H38O4	1000377-72-3	80
5			1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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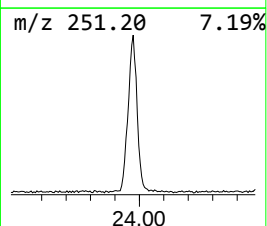
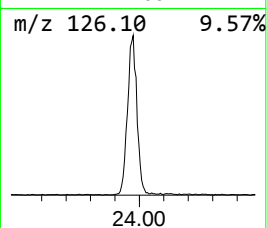
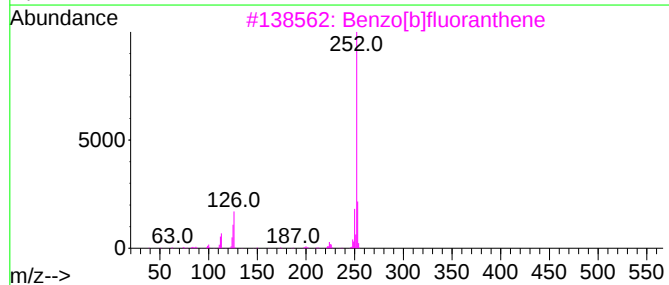
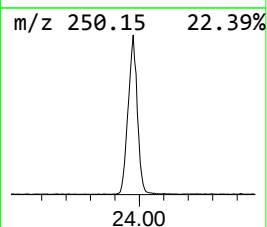
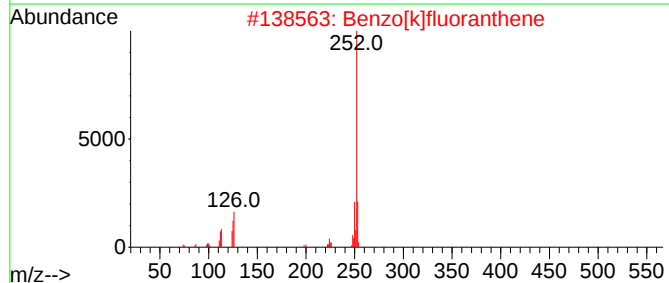
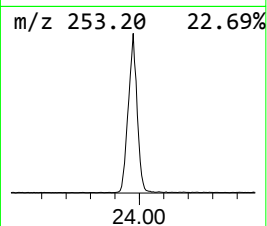
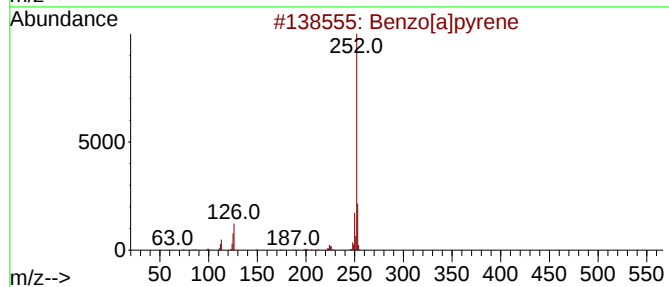
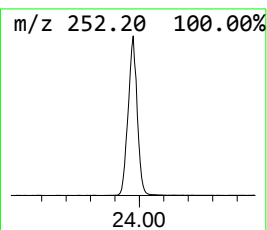
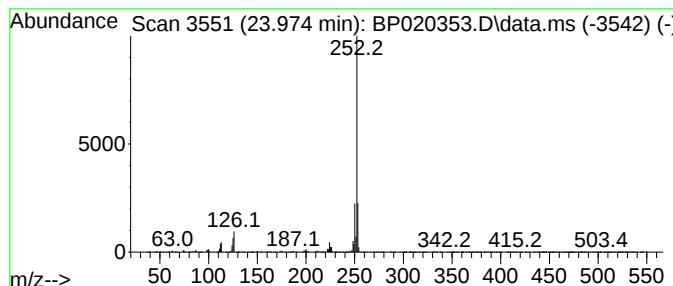
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Benzo[a]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.974	33.09 ng/ul	2381240	Perylene-d12	24.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[a]pyrene	252	C20H12	000050-32-8	96
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5			Benzo[e]pyrene	252	C20H12	000192-97-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
Data File : BP020353.D
Acq On : 13 May 2024 12:47
Operator : MA/JU
Sample : P2372-23DL 5X
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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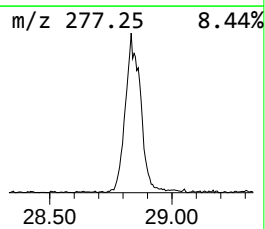
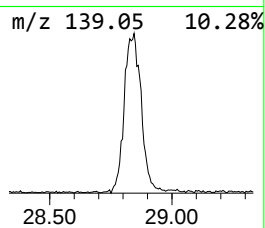
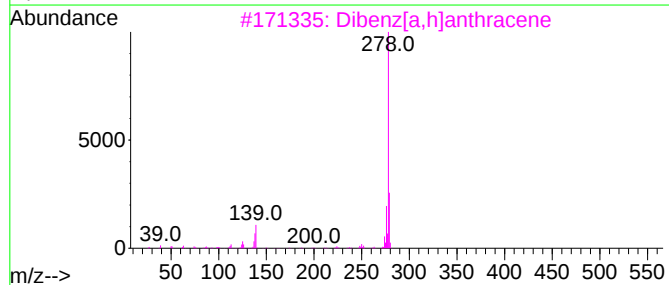
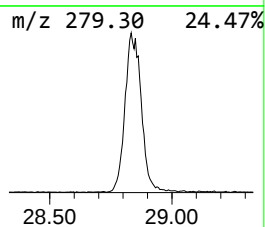
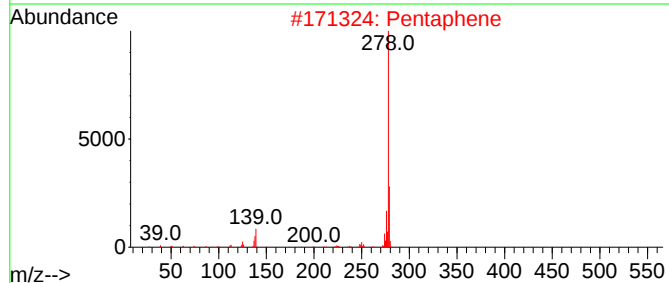
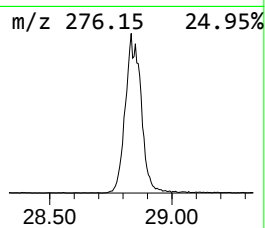
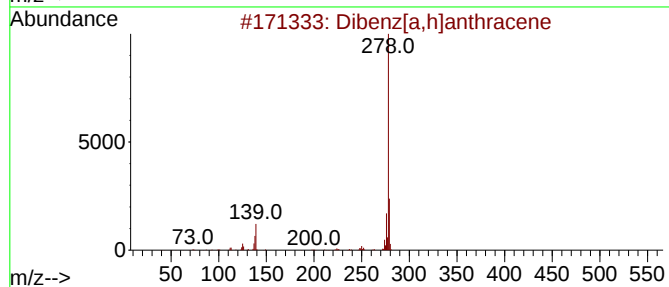
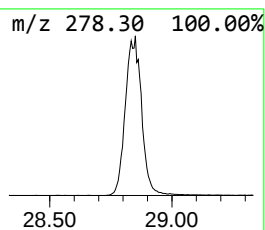
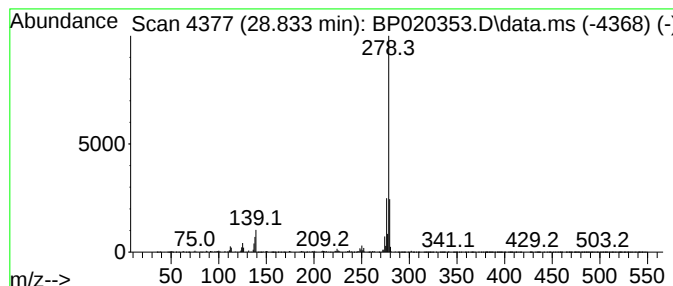
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 23 Dibenz[a,h]anthracene Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.833	8.88 ng/ul	639228	Perylene-d12	24.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenz[a,h]anthracene	278	C22H14	000053-70-3	98
2			Pentaphene	278	C22H14	000222-93-5	98
3			Dibenz[a,h]anthracene	278	C22H14	000053-70-3	98
4			Picene	278	C22H14	000213-46-7	97
5			Benzo[b]chrysene	278	C22H14	000214-17-5	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051324\
 Data File : BP020353.D
 Acq On : 13 May 2024 12:47
 Operator : MA/JU
 Sample : P2372-23DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43Y3DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Bis(2-chloroeth...	7.063	4.1	ng/ul	106097	1	7.616	518681	20.0
Benzene, 1,3-di...	7.499	6.9	ng/ul	179321	1	7.616	518681	20.0
Benzene, 1,4-di...	7.958	19.2	ng/ul	497438	1	7.616	518681	20.0
1-Propanamine, ...	8.422	15.1	ng/ul	391542	1	7.616	518681	20.0
Ethane, hexachl...	8.663	16.6	ng/ul	429692	1	7.616	518681	20.0
Methane, bis(2-...	9.822	5.7	ng/ul	211297	2	10.381	743254	20.0
Naphthalene	10.434	26.5	ng/ul	984008	2	10.381	743254	20.0
Phenol, 4-chlor...	11.704	15.6	ng/ul	579253	2	10.381	743254	20.0
Phenol, 2,4,6-t...	12.693	7.8	ng/ul	396976	3	14.281	1020200	20.0
Phenol, 2,4,5-t...	12.775	16.7	ng/ul	851472	3	14.281	1020200	20.0
Naphthalene, 1-...	13.134	9.4	ng/ul	480572	3	14.281	1020200	20.0
Benzene, 2-meth...	13.881	21.1	ng/ul	1078660	3	14.281	1020200	20.0
Acena phthene	14.345	22.0	ng/ul	1122950	3	14.281	1020200	20.0
unknown-01	14.510	27.1	ng/ul	1381010	3	14.281	1020200	20.0
Benzene, 1-meth...	14.704	13.3	ng/ul	677849	3	14.281	1020200	20.0
Diethyl Phthalate	15.145	30.8	ng/ul	1568320	3	14.281	1020200	20.0
Benzene, 1-chlo...	15.363	34.1	ng/ul	1738040	3	14.281	1020200	20.0
Benzene, hexach...	16.387	32.1	ng/ul	2050300	4	17.116	1277920	20.0
1,2-Benzenedica...	18.145	26.9	ng/ul	1721380	4	17.116	1277920	20.0
Fluoranthene	19.275	36.2	ng/ul	2310220	4	17.116	1277920	20.0
Bis(2-ethylhexy...	21.533	16.6	ng/ul	2585070	5	21.616	3119030	20.0
Benzo[a]pyrene	23.974	33.1	ng/ul	2381240	6	24.963	1439440	20.0
Dibenz[a,h]anth...	28.833	8.9	ng/ul	639228	6	24.963	1439440	20.0