

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
 Data File : BN031258.D
 Acq On : 27 Apr 2024 00:24
 Operator : MA/JU
 Sample : P2161-20
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 CORD7

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_N\Methods\SFAM-EPA-BN041924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN031258.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.552	92	96	113	rBV	365664	681247	0.38%	0.078%
2	5.058	329	352	356	rBV4	29024973	166907703	92.61%	19.215%
3	6.316	559	566	575	rBV	180039	299766	0.17%	0.035%
4	6.611	610	616	629	rBV	223992	412606	0.23%	0.048%
5	6.722	629	635	643	rVV	101631	180525	0.10%	0.021%
6	6.805	643	649	663	rVB	361595	665867	0.37%	0.077%
7	6.993	670	681	692	rBV	938585	2289559	1.27%	0.264%
8	7.169	702	711	722	rBV	1085628	2311020	1.28%	0.266%
9	7.546	761	775	786	rBV3	18119066	71881636	39.88%	8.275%
10	7.758	786	811	815	rVV2	29003058	156459994	86.81%	18.012%
11	8.093	842	868	870	rBV2	29796419	180229469	100.00%	20.748%
12	8.340	903	910	927	rVB	1072934	1732210	0.96%	0.199%
13	8.793	979	987	998	rBV	1706614	2769575	1.54%	0.319%
14	9.116	1034	1042	1048	rBV	977203	1559514	0.87%	0.180%
15	9.340	1074	1080	1083	rBV	135032	214127	0.12%	0.025%
16	9.381	1083	1087	1090	rVV2	168449	311827	0.17%	0.036%
17	9.422	1090	1094	1101	rVV	514064	911572	0.51%	0.105%
18	9.505	1101	1108	1113	rVV	1367995	2318527	1.29%	0.267%
19	9.552	1113	1116	1126	rVB	225004	358127	0.20%	0.041%
20	10.040	1191	1199	1208	rBV	1667010	3325670	1.85%	0.383%
21	10.216	1221	1229	1232	rBV	179603	303054	0.17%	0.035%
22	10.346	1232	1251	1255	rVB2	28466411	122274421	67.84%	14.077%
23	10.440	1260	1267	1271	rVV	1466852	2210382	1.23%	0.254%
24	10.481	1271	1274	1279	rVV	130310	182313	0.10%	0.021%
25	10.563	1279	1288	1299	rVB	1664841	3107072	1.72%	0.358%
26	10.828	1319	1333	1338	rBV2	20337902	58169565	32.28%	6.697%
27	12.446	1592	1608	1617	rBV	4655423	7834133	4.35%	0.902%
28	12.693	1644	1650	1661	rBV2	138032	283993	0.16%	0.033%
29	13.063	1705	1713	1725	rBV	8189654	13482289	7.48%	1.552%
30	13.693	1808	1820	1834	rBV	3564360	5206667	2.89%	0.599%
31	13.963	1853	1866	1875	rBV	4315688	6476259	3.59%	0.746%
32	14.269	1910	1918	1932	rVB	2029230	3019952	1.68%	0.348%
33	14.487	1949	1955	1966	rBV	388054	642761	0.36%	0.074%
34	14.587	1966	1972	1977	rVB	196722	282987	0.16%	0.033%
35	14.763	1996	2002	2017	rBV3	179364	389008	0.22%	0.045%
36	15.269	2081	2088	2103	rVB	5233239	7751935	4.30%	0.892%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

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37	15.392	2103	2109	2124	rBV	1820037	2521507	1.40%	0.290%
38	17.022	2376	2386	2392	rBV	2584516	3558018	1.97%	0.410%
39	17.122	2396	2403	2413	rVB2	5686633	8086711	4.49%	0.931%
40	17.922	2534	2539	2547	rBV	220381	329195	0.18%	0.038%
41	19.210	2754	2758	2766	rBV2	122100	187403	0.10%	0.022%
42	19.428	2788	2795	2801	rBV	7122025	9594547	5.32%	1.105%
43	20.957	3051	3055	3064	rBV4	150826	279050	0.15%	0.032%
44	21.257	3100	3106	3117	rBV2	2655879	3765315	2.09%	0.433%
45	23.321	3454	3457	3466	rVB	120486	206508	0.11%	0.024%
46	23.963	3555	3566	3581	rVB2	3684320	8964399	4.97%	1.032%
47	24.151	3590	3598	3611	rVB	1571906	3710906	2.06%	0.427%

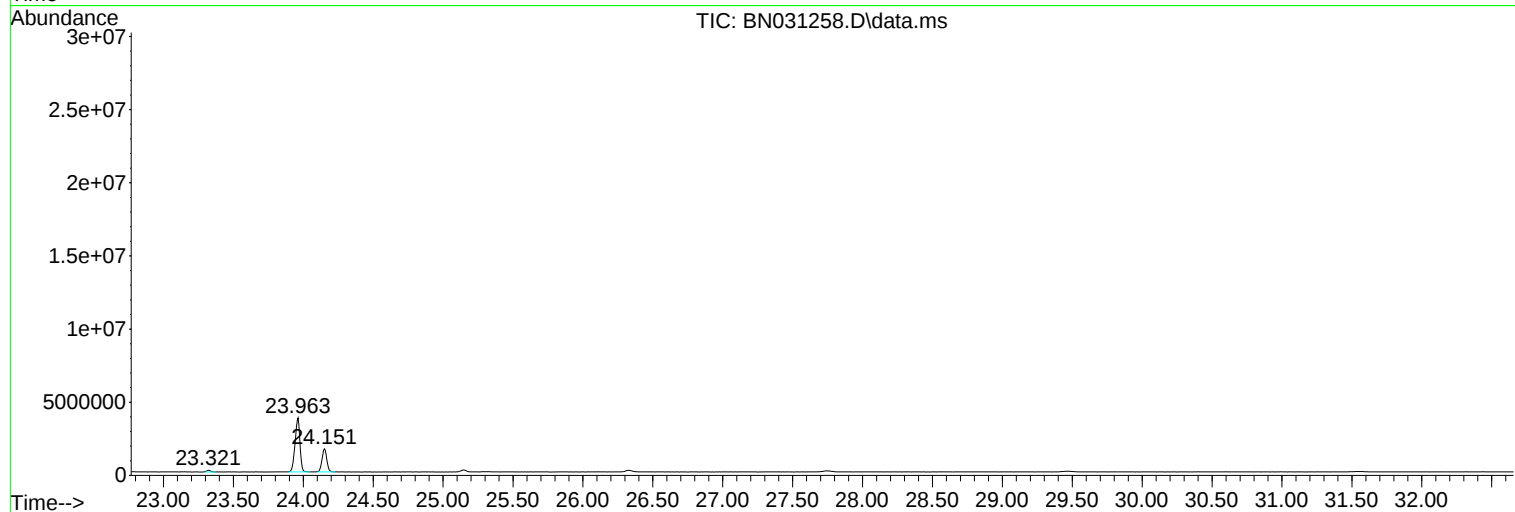
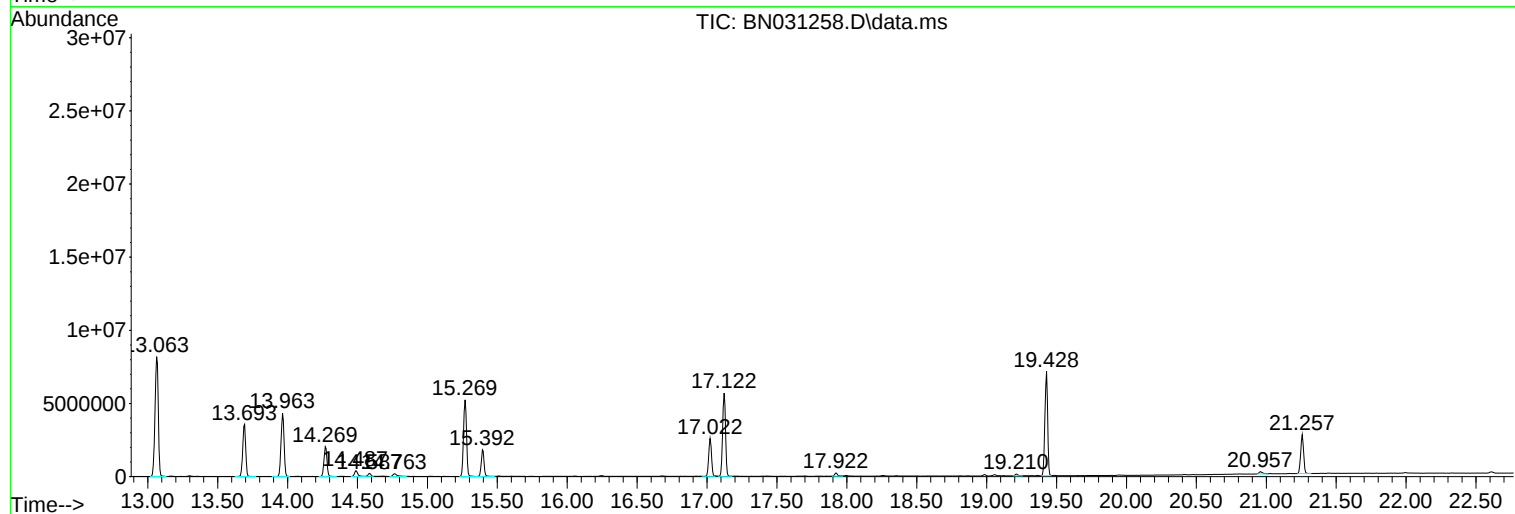
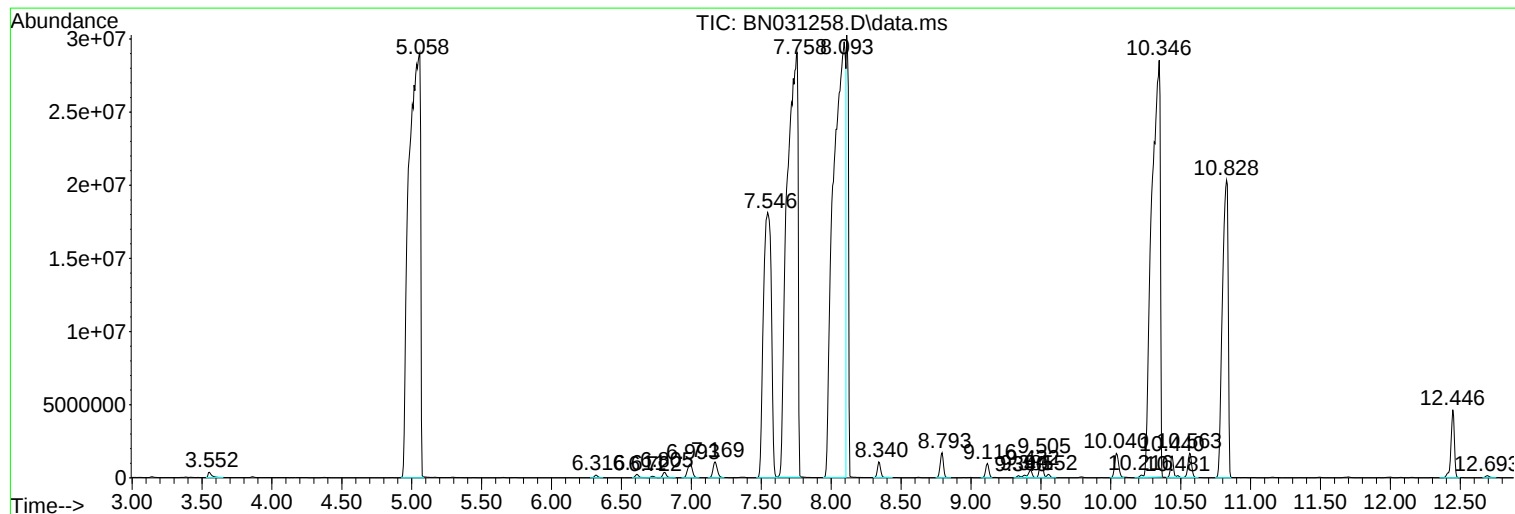
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : MA/JU
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
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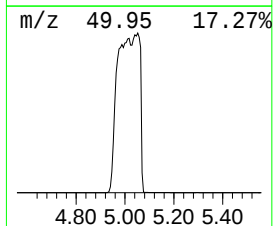
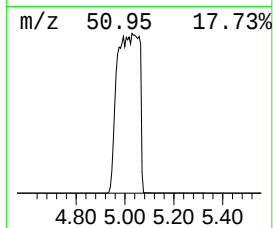
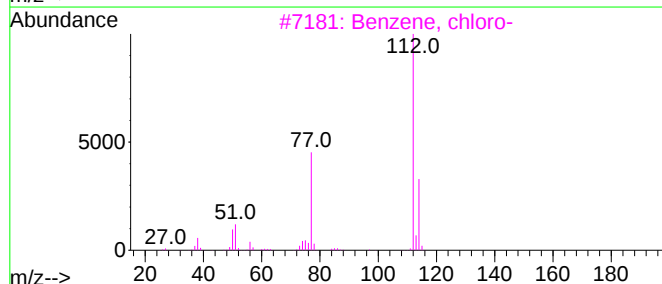
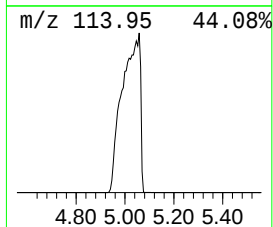
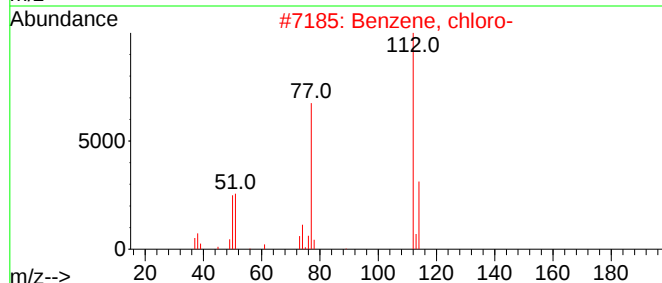
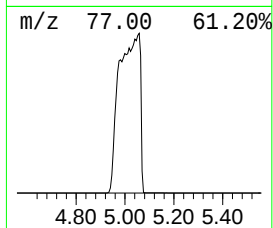
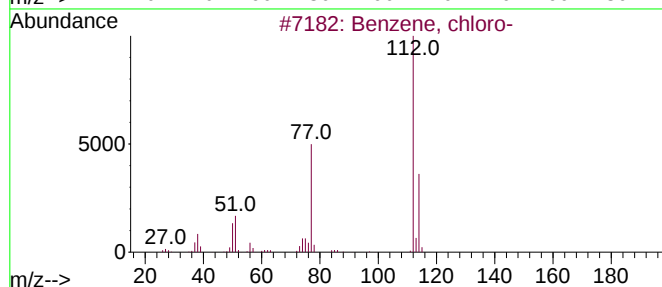
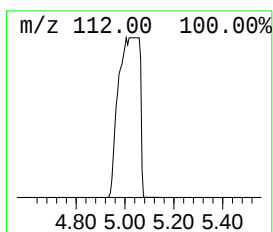
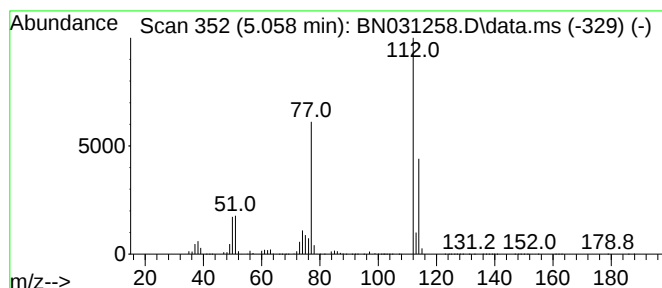
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.058	21.34 ng/ul	166908000	1,4-Dichlorobenzene-d4	7.652

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	93
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	90



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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

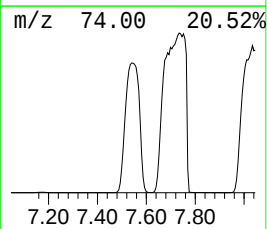
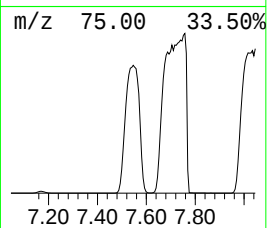
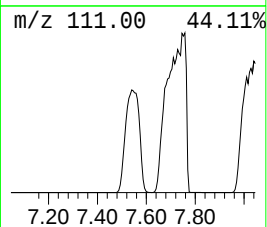
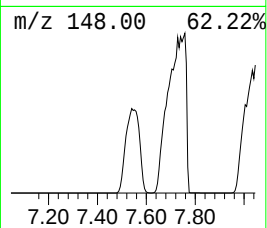
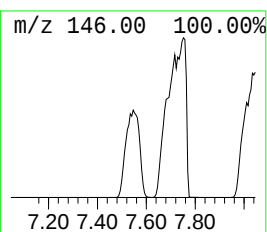
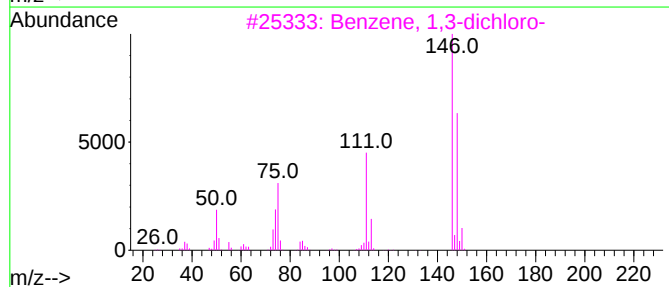
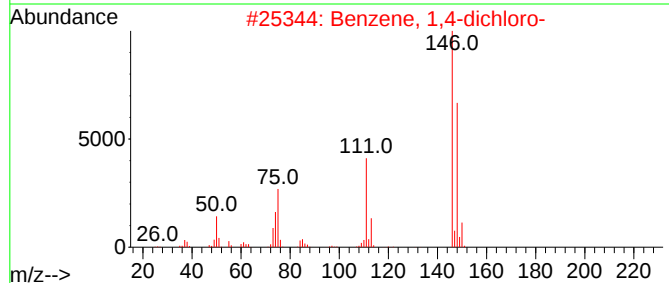
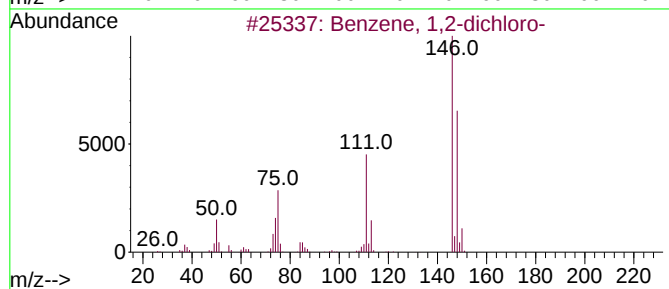
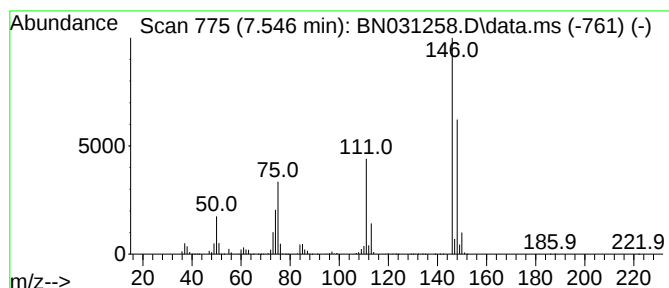
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.546	9.19 ng/ul	71881600	1,4-Dichlorobenzene-d4	7.652

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

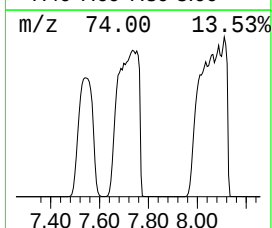
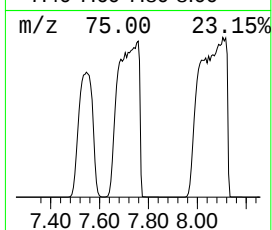
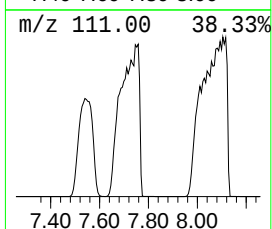
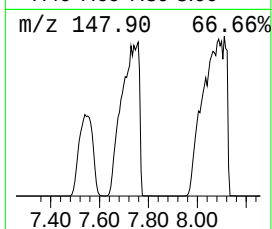
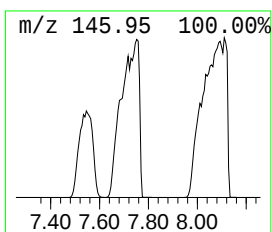
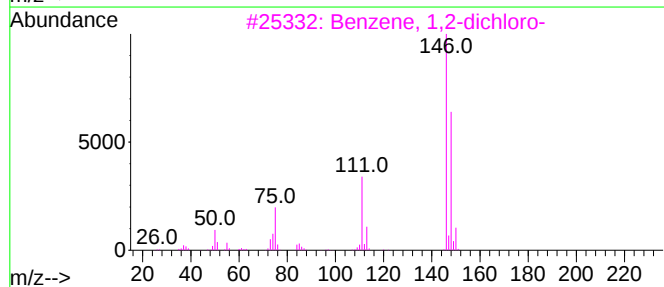
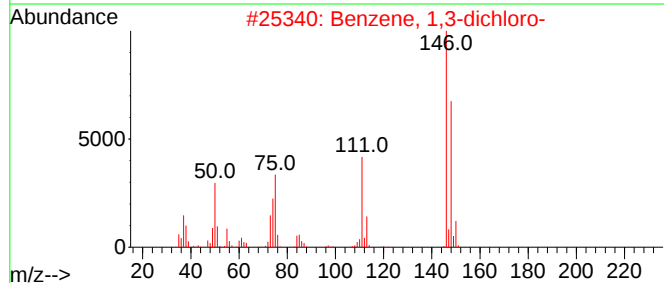
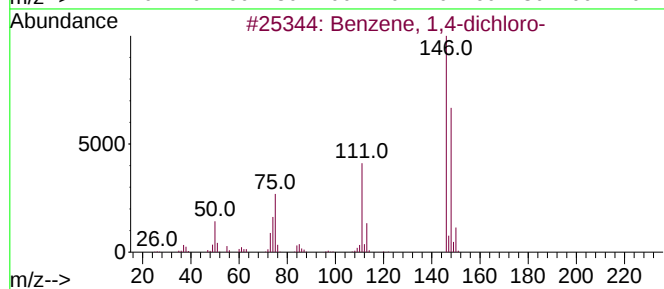
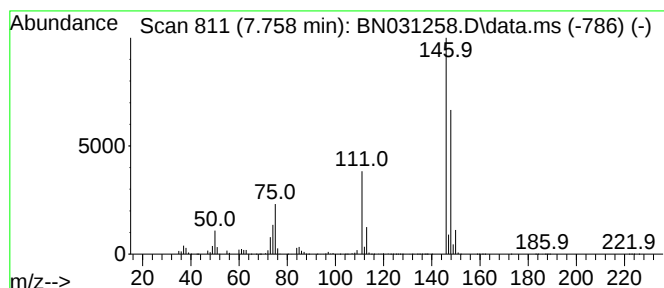
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.758	20.00 ng/ul	156460000	1,4-Dichlorobenzene-d4	7.652

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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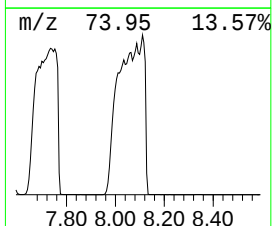
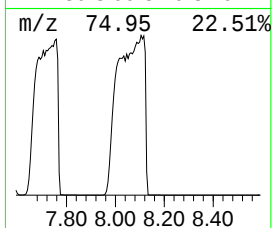
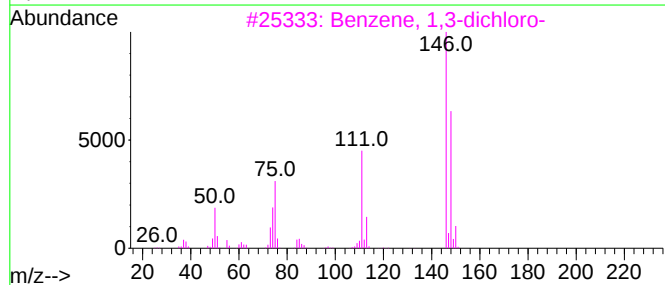
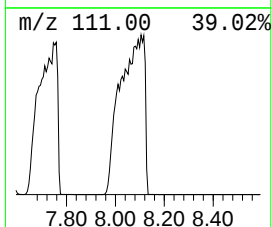
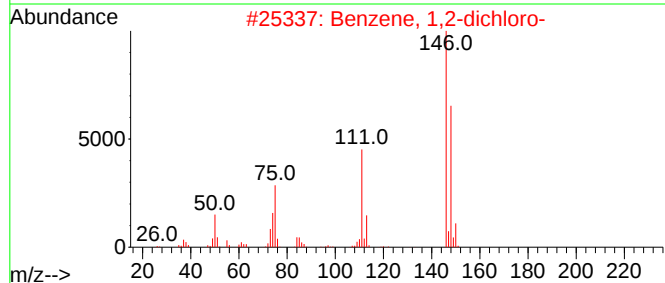
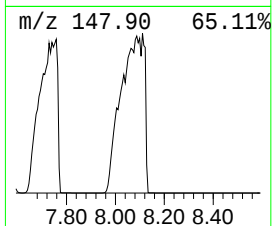
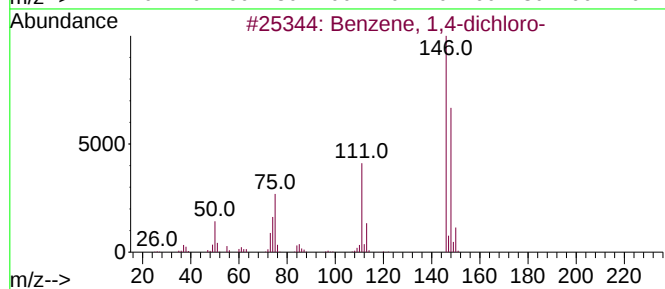
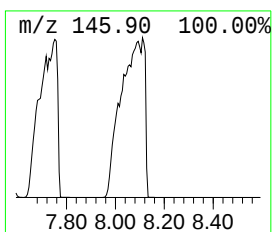
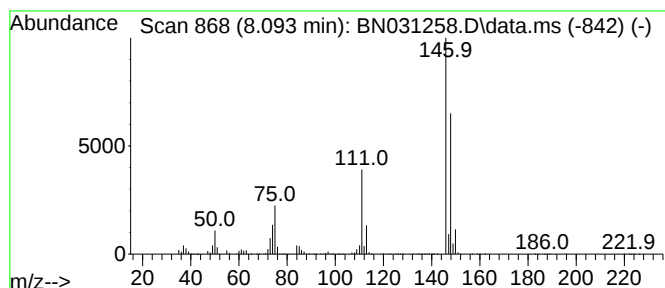
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.093	23.04 ng/ul	180229000	1,4-Dichlorobenzene-d4	7.652

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	97
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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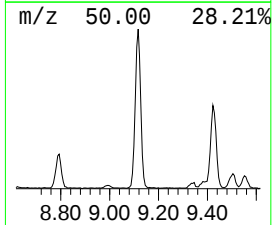
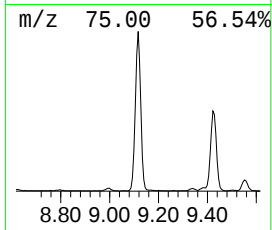
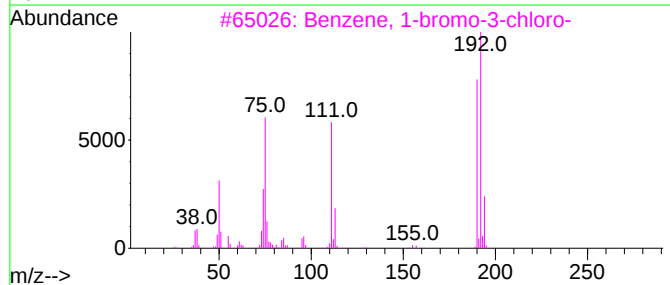
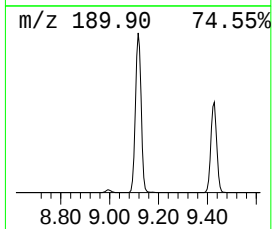
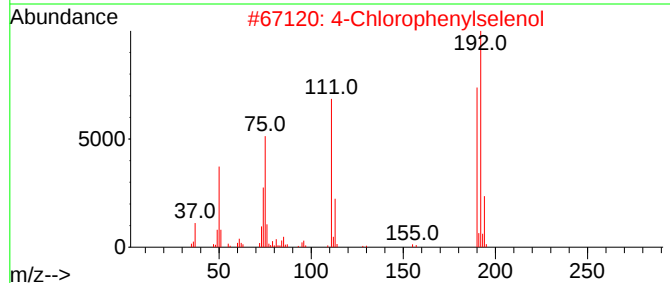
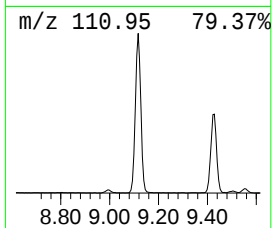
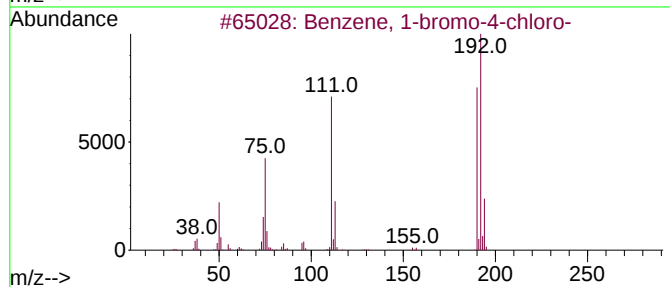
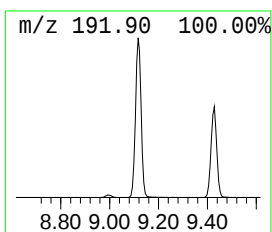
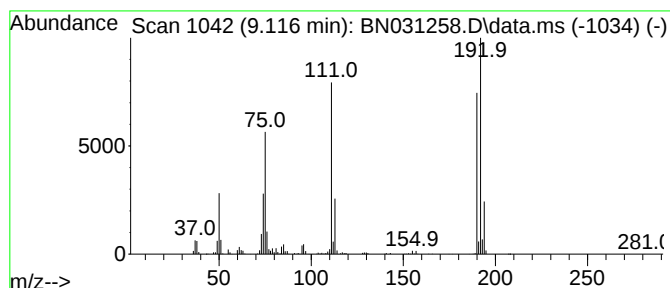
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.116	14.11 ng/ul	1559510	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
2			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
3			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
Acq On : 27 Apr 2024 00:24
Operator : MA/JU
Sample : P2161-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORD7

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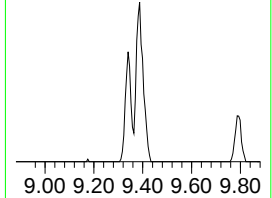
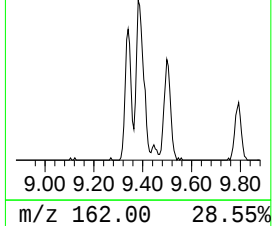
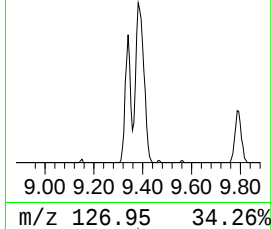
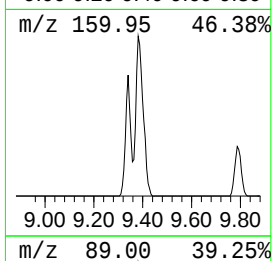
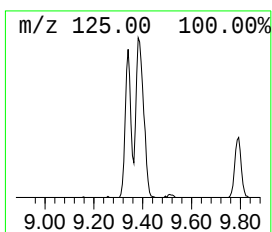
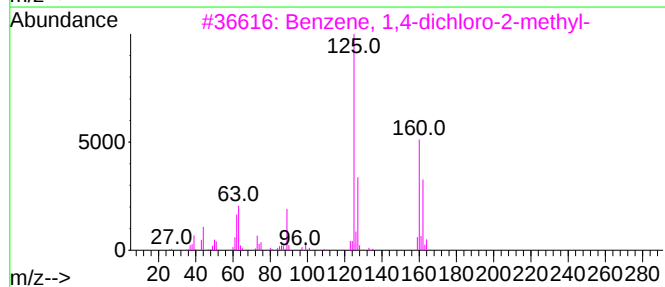
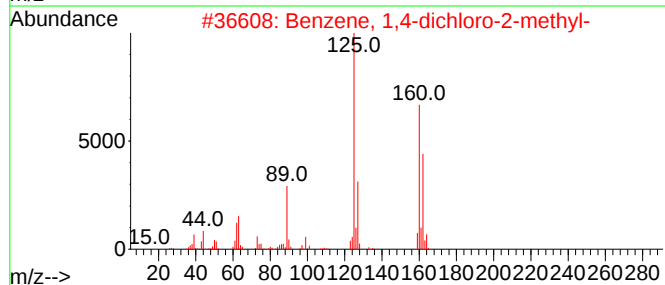
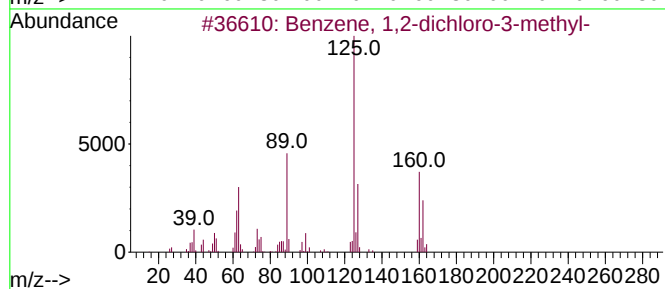
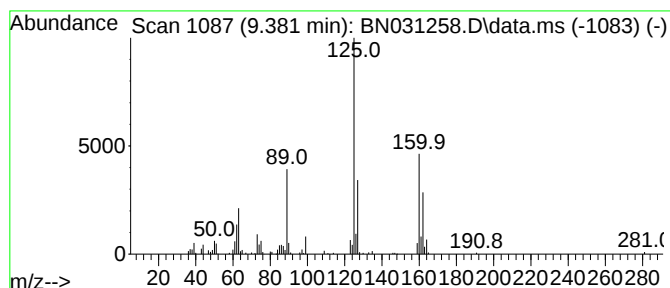
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TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1,2-dichloro-3-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.381	2.82 ng/ul	311827	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96
2			Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
3			Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
4			Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	95
5			Benzene, 1-chloro-2-(chloromethyl)-	160	C7H6Cl2	000611-19-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
Acq On : 27 Apr 2024 00:24
Operator : MA/JU
Sample : P2161-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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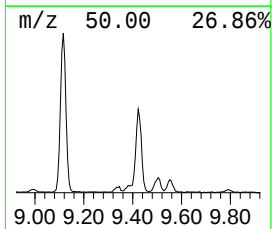
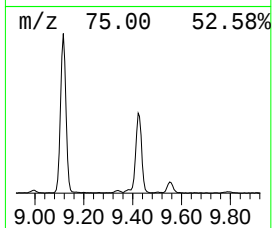
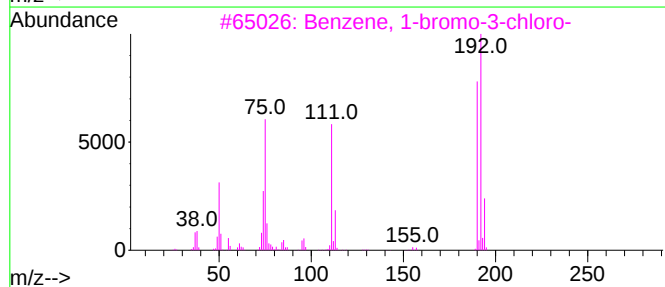
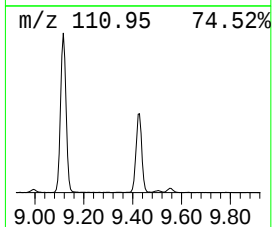
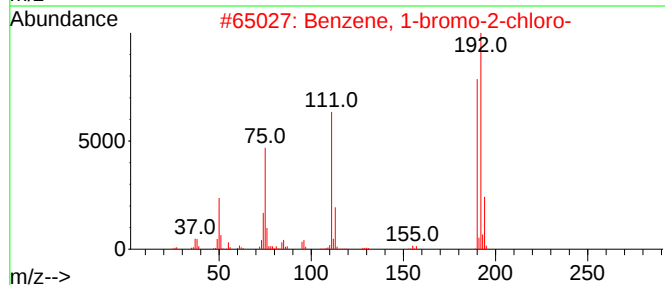
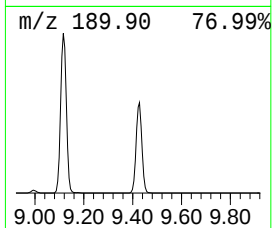
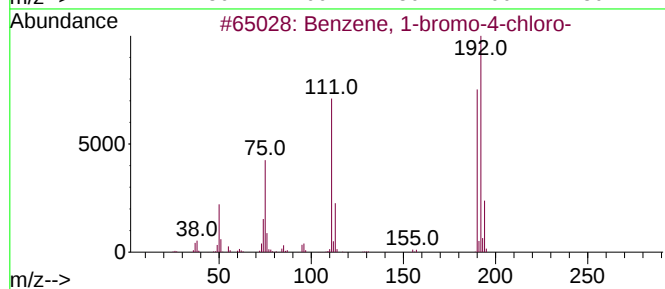
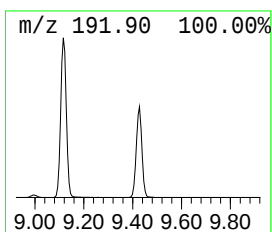
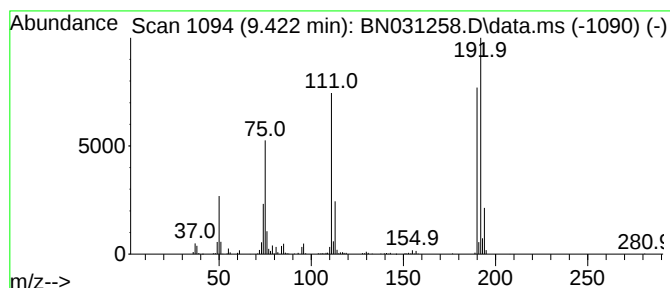
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1-bromo-2-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.422	8.25 ng/ul	911572	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
2			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
4			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	96
5			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
Acq On : 27 Apr 2024 00:24
Operator : MA/JU
Sample : P2161-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORD7

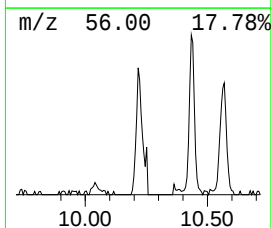
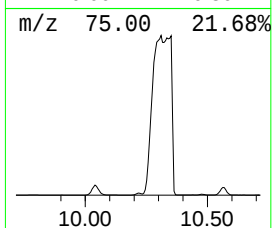
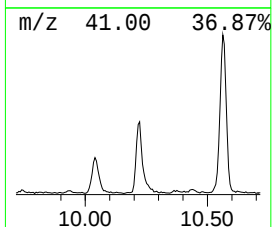
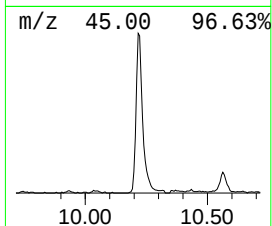
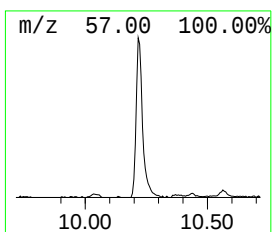
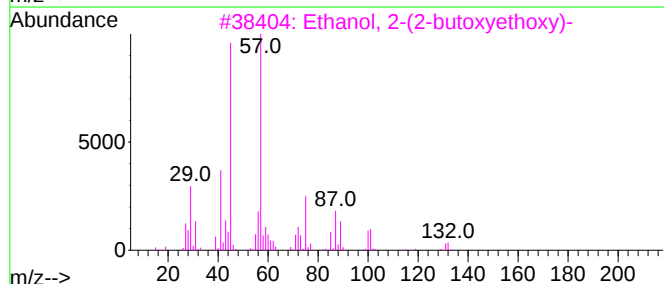
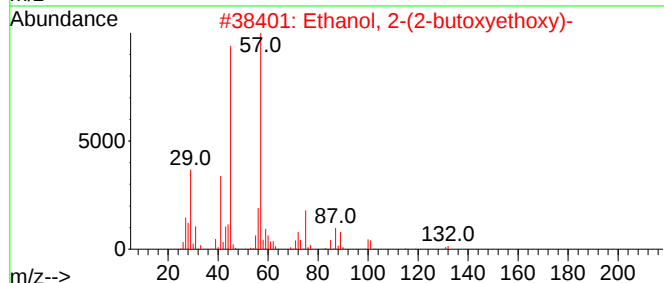
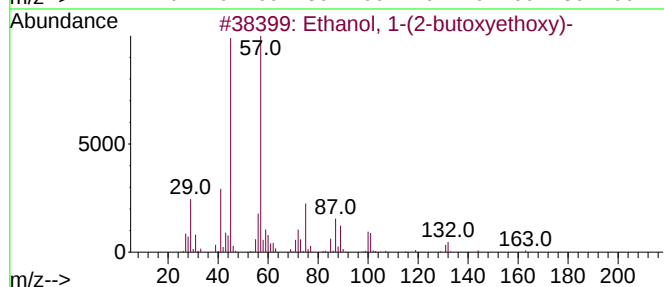
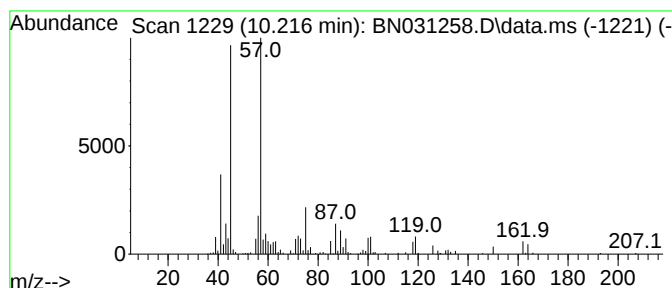
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.216	2.74 ng/ul	303054	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	80
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	80
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	80
4			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	80
5			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
Acq On : 27 Apr 2024 00:24
Operator : MA/JU
Sample : P2161-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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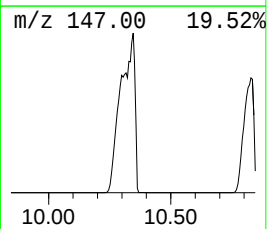
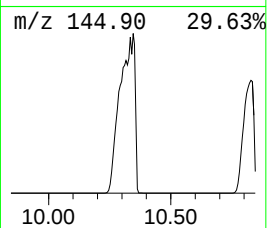
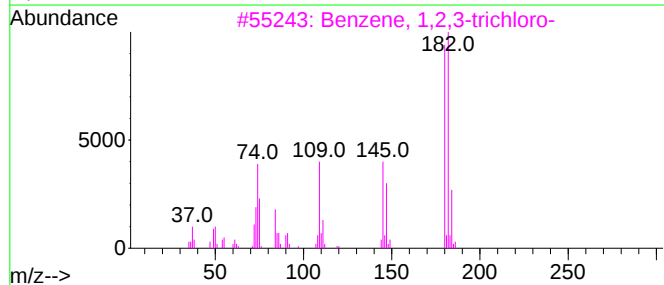
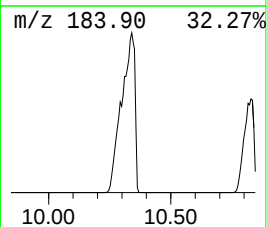
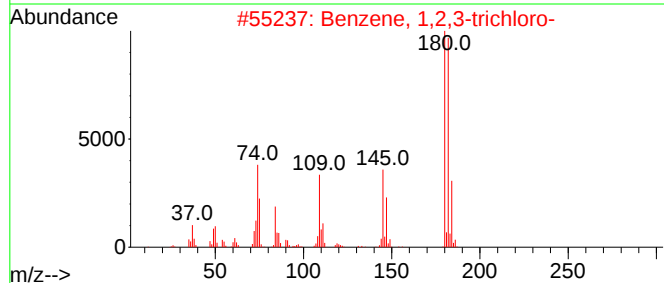
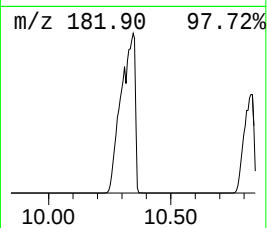
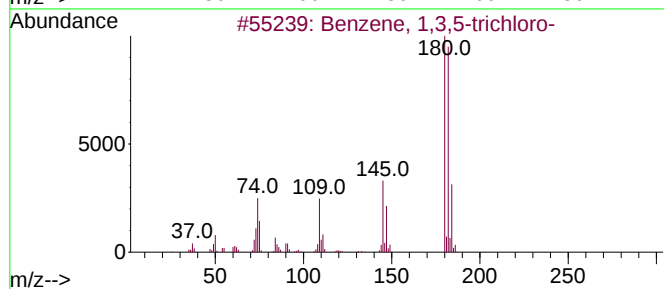
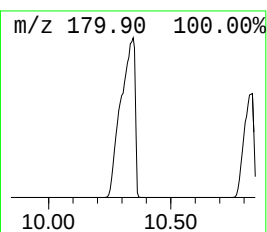
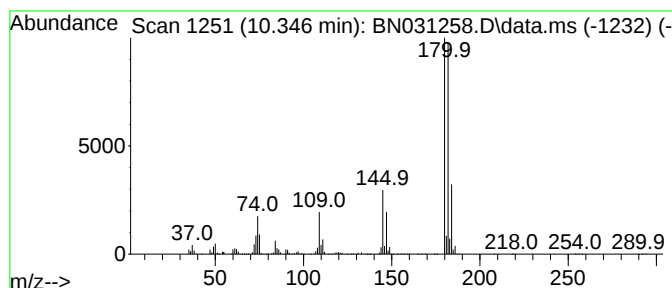
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.346	1106.36 ng/ul	122274000	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
Acq On : 27 Apr 2024 00:24
Operator : MA/JU
Sample : P2161-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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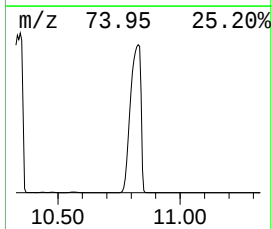
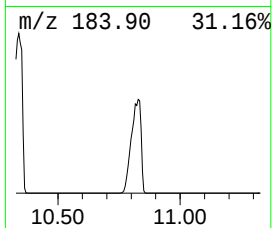
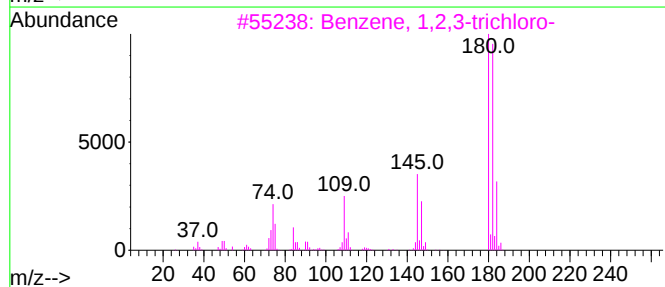
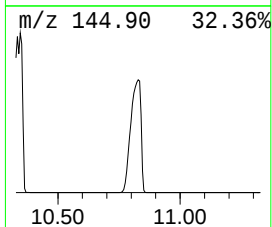
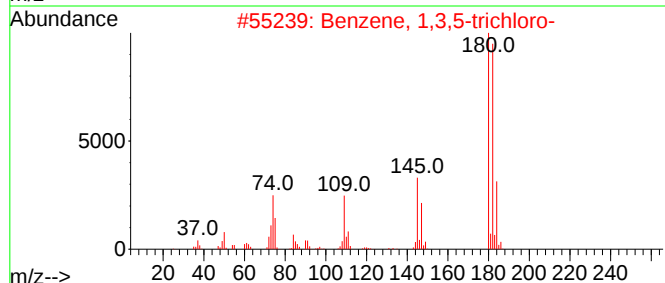
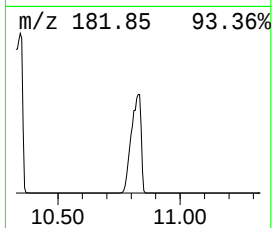
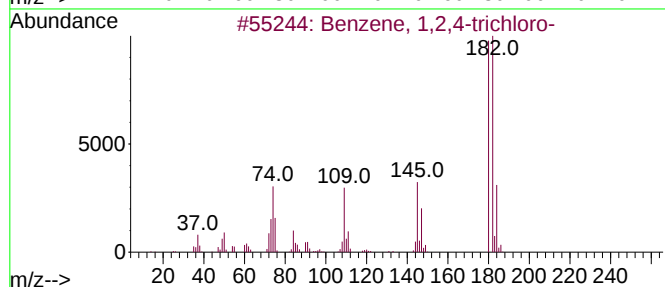
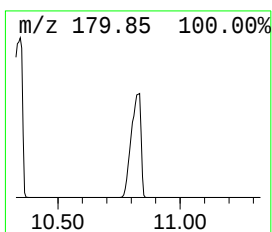
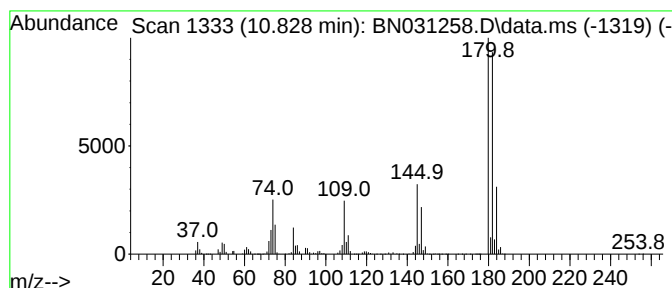
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TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1,2,4-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.828	526.33 ng/ul	58169600	Naphthalene-d8	10.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	98
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
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ClientSampleId :
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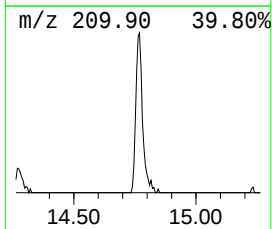
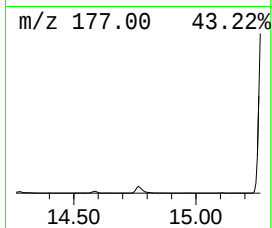
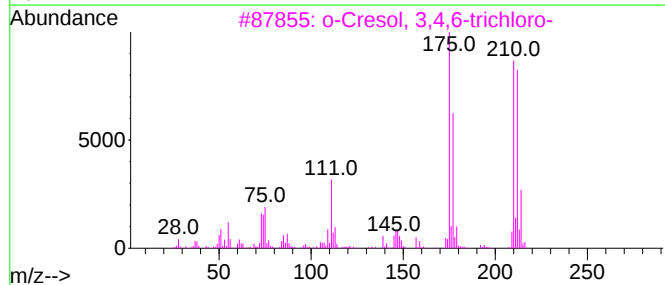
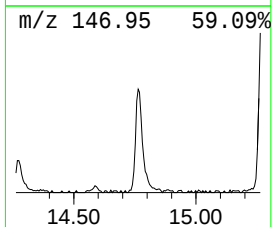
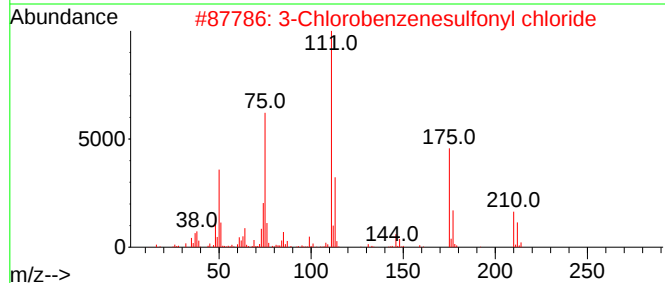
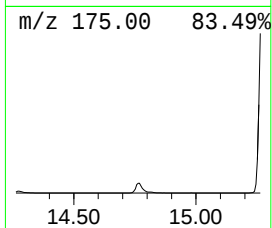
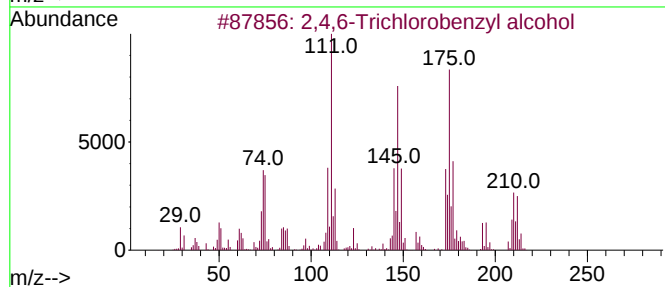
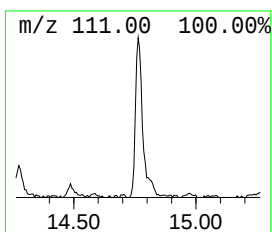
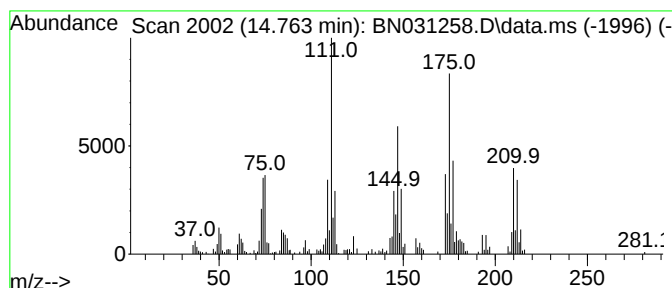
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2,4,6-Trichlorobenzyl alcohol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.763	2.58 ng/ul	389008	Acenaphthene-d10	14.269

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,6-Trichlorobenzyl alcohol	210	C7H5Cl3O	217479-60-2	81	
2	3-Chlorobenzenesulfonyl chloride	210	C6H4Cl2O2S	002888-06-4	64	
3	o-Cresol, 3,4,6-trichloro-	210	C7H5Cl3O	000643-14-1	53	
4	2,5-Dichloro-.alpha.-methylbenzy...	190	C8H8Cl2O	001475-12-3	53	
5	Benzenemethanol, 2,3-dichloro-.a...	190	C8H8Cl2O	054798-91-3	53	



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042624\
Data File : BN031258.D
Acq On : 27 Apr 2024 00:24
Operator : MA/JU
Sample : P2161-20
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
CORD7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.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
Benzene, chloro-	5.058	21.3	ng/ul	166908000	1	7.652	156460000	20.0
Benzene, 1,2-di...	7.546	9.2	ng/ul	71881600	1	7.652	156460000	20.0
Benzene, 1,4-di...	7.758	20.0	ng/ul	156460000	1	7.652	156460000	20.0
Benzene, 1,3-di...	8.093	23.0	ng/ul	180229000	1	7.652	156460000	20.0
Benzene, 1-brom...	9.116	14.1	ng/ul	1559510	2	10.434	2210380	20.0
Benzene, 1,2-di...	9.381	2.8	ng/ul	311827	2	10.434	2210380	20.0
Benzene, 1-brom...	9.422	8.3	ng/ul	911572	2	10.434	2210380	20.0
Ethanol, 1-(2-b...	10.216	2.7	ng/ul	303054	2	10.434	2210380	20.0
Benzene, 1,3,5-...	10.346	1106.4	ng/ul	122274000	2	10.434	2210380	20.0
Benzene, 1,2,4-...	10.828	526.3	ng/ul	58169600	2	10.434	2210380	20.0
2,4,6-Trichloro...	14.763	2.6	ng/ul	389008	3	14.269	3019950	20.0