

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
 Data File : BM034612.D
 Acq On : 11 Apr 2022 16:27
 Operator : CG/JU
 Sample : N2338-02
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0J71

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_M\Methods\SFAM-EPA-BM040922.M
 Title : SVOA CALIBRATION

Signal : TIC: BM034612.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.690	108	111	128	rBV	140607	251797	0.11%	0.026%
2	5.196	355	367	371	rBV3	41222325	137457468	59.55%	14.194%
3	6.554	591	598	607	rBV	261726	433010	0.19%	0.045%
4	7.078	676	687	697	rBV2	309301	786096	0.34%	0.081%
5	7.219	701	711	728	rBV	5656815	11291002	4.89%	1.166%
6	7.419	738	745	758	rVB	510586	1019927	0.44%	0.105%
7	7.795	797	809	820	rBV	31876784	80966140	35.07%	8.361%
8	7.984	820	841	847	rBV6	43012406	208170693	90.18%	21.496%
9	8.319	878	898	904	rBV5	43231376	230838441	100.00%	23.836%
10	8.601	940	946	960	rVB	398717	731082	0.32%	0.075%
11	9.148	1015	1039	1043	rBV	24021738	69181195	29.97%	7.144%
12	9.307	1062	1066	1074	rVB	188930	301058	0.13%	0.031%
13	9.390	1074	1080	1089	rVB	399449	658500	0.29%	0.068%
14	9.719	1128	1136	1142	rVV	3167670	5492985	2.38%	0.567%
15	9.778	1142	1146	1153	rVV	580710	957344	0.41%	0.099%
16	10.319	1231	1238	1246	rBV	736912	1303395	0.56%	0.135%
17	10.595	1271	1285	1290	rBV	40344459	104633933	45.33%	10.805%
18	10.707	1298	1304	1310	rBV	656537	995725	0.43%	0.103%
19	10.819	1314	1323	1346	rVB	8285216	16742109	7.25%	1.729%
20	11.101	1360	1371	1384	rBV	24618966	45146225	19.56%	4.662%
21	11.307	1396	1406	1428	rBV	10276019	18360428	7.95%	1.896%
22	11.513	1430	1441	1460	rVB	1912478	3428074	1.49%	0.354%
23	12.048	1526	1532	1543	rBV	396367	645241	0.28%	0.067%
24	12.725	1633	1647	1670	rBV2	1458547	2706947	1.17%	0.280%
25	12.960	1681	1687	1696	rBV	244747	409084	0.18%	0.042%
26	13.330	1743	1750	1760	rBV	2015252	3036166	1.32%	0.314%
27	13.936	1847	1853	1867	rVB	1324397	1907769	0.83%	0.197%
28	14.219	1894	1901	1914	rBV	1635263	2324933	1.01%	0.240%
29	14.524	1946	1953	1963	rBV	862113	1269238	0.55%	0.131%
30	15.519	2115	2122	2135	rBV	1840738	2772294	1.20%	0.286%
31	15.630	2135	2141	2155	rVV	507918	857388	0.37%	0.089%
32	17.265	2412	2419	2427	rBV	963513	1350427	0.59%	0.139%
33	17.365	2430	2436	2450	rBV	1944319	2891942	1.25%	0.299%
34	19.654	2818	2825	2844	rBV	2359098	3431646	1.49%	0.354%
35	21.442	3125	3129	3134	rBV	762002	1129723	0.49%	0.117%
36	23.671	3501	3508	3525	rBV	1355724	3052452	1.32%	0.315%

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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_M\Methods\SFAM-EPA-BM040922.M
Title : SVOA CALIBRATION

37	23.824	3528	3534	3549	rVB2	654494	1494825	0.65%	0.154%
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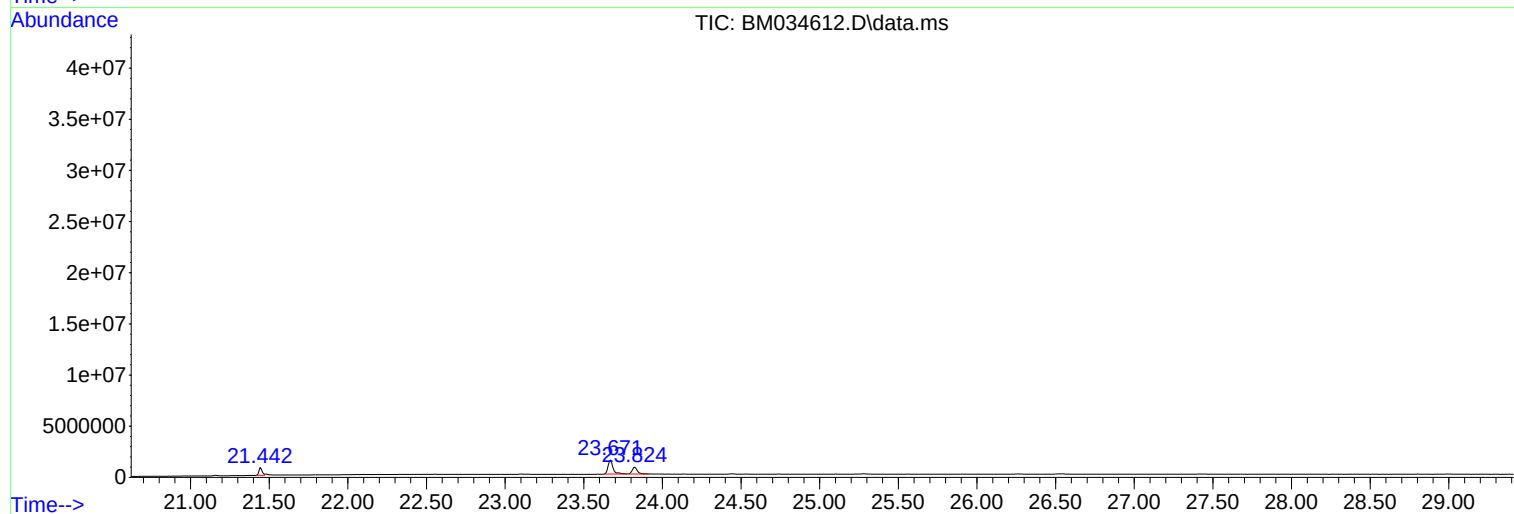
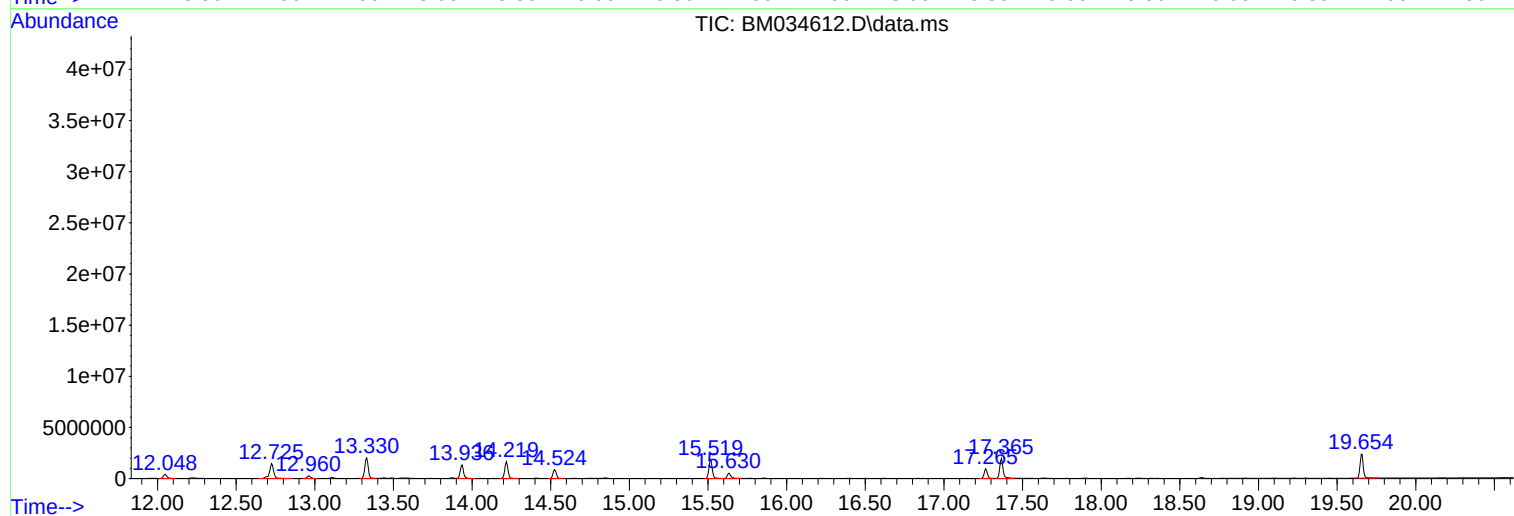
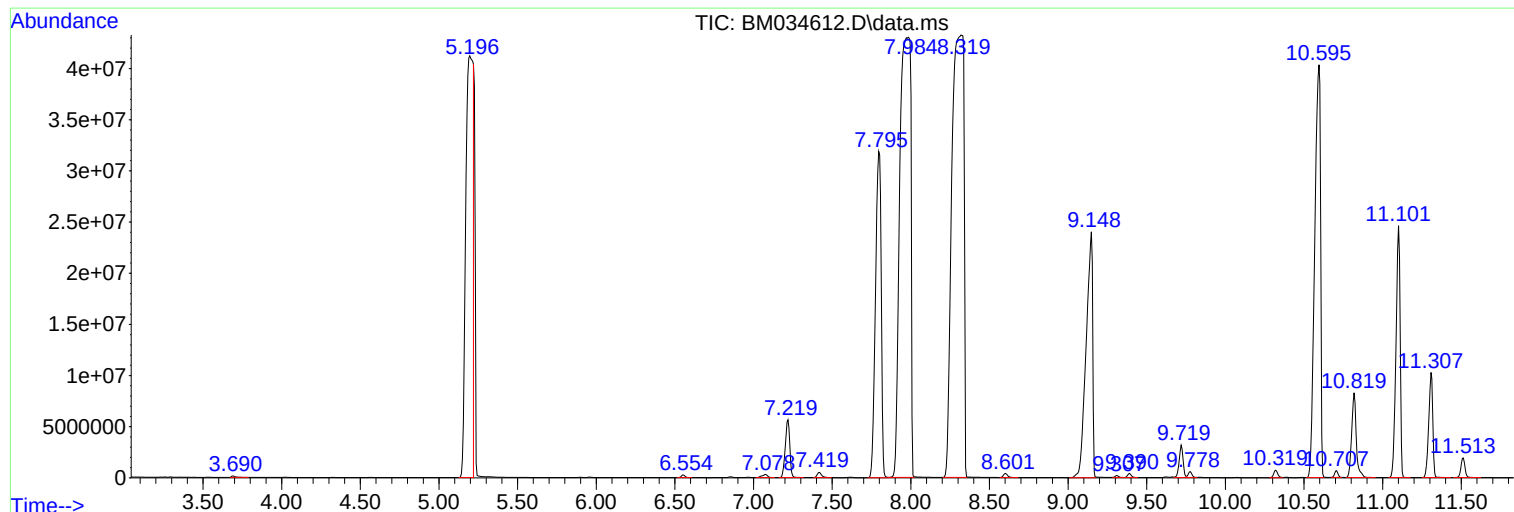
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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



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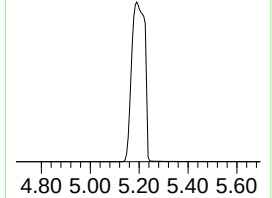
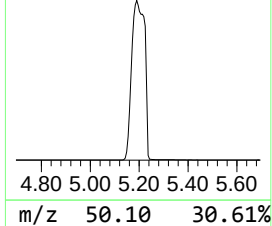
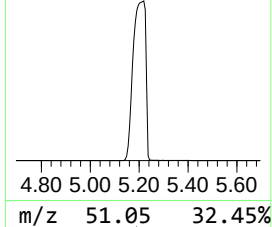
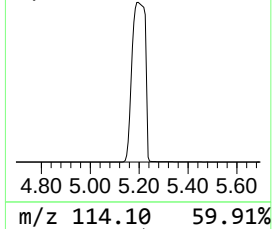
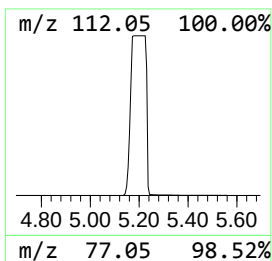
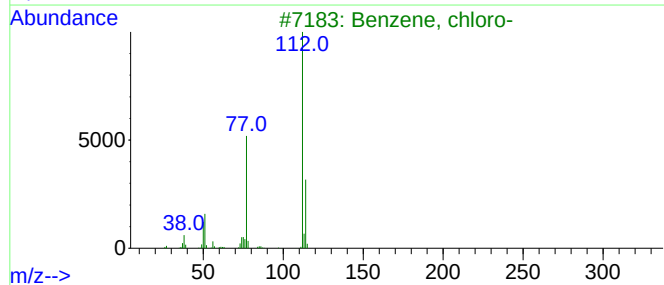
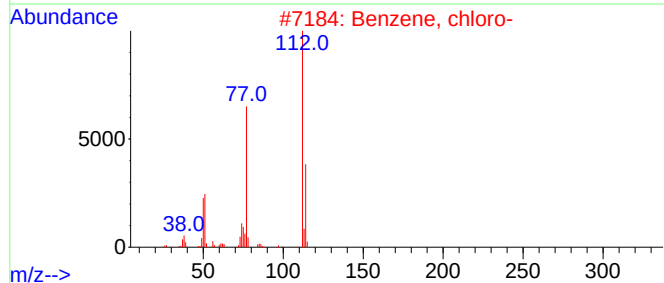
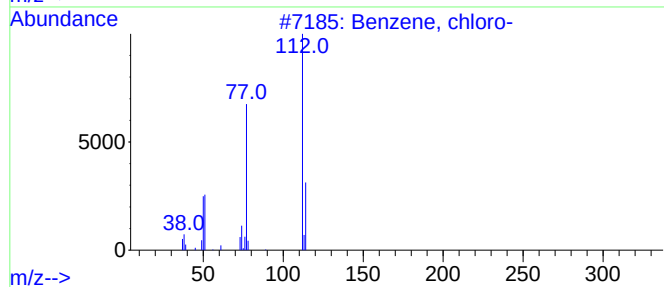
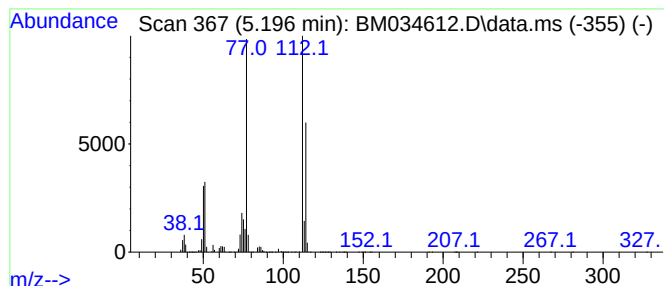
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.196	13.21 ng/ul	137457000	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	93
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	93
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
5			Benzene, chloro-	112	C6H5Cl	000108-90-7	87



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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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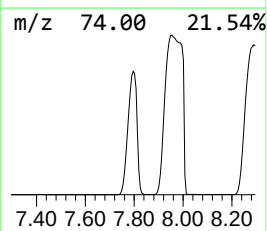
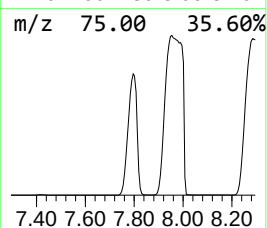
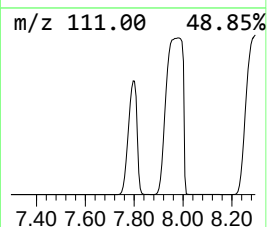
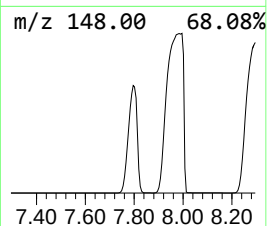
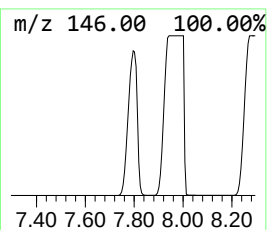
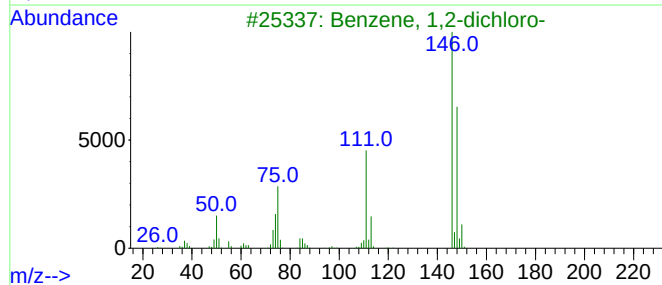
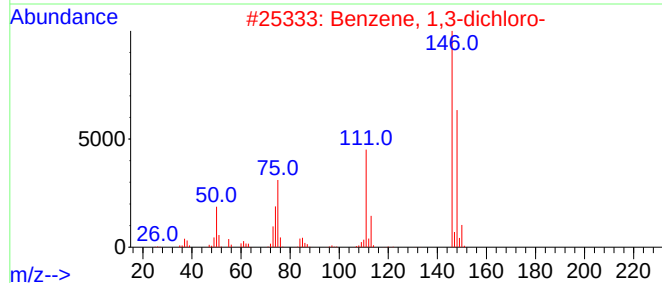
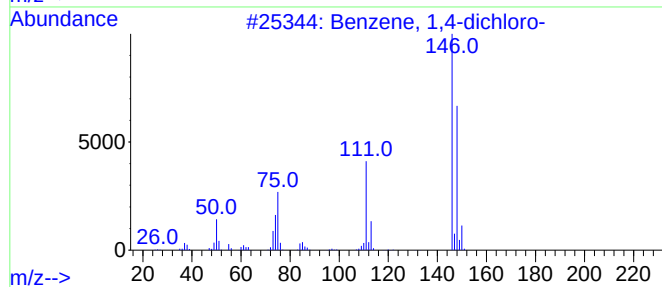
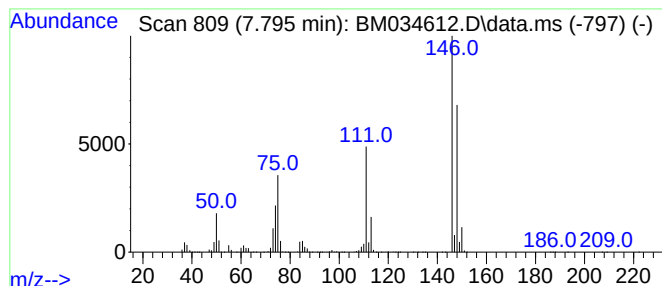
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.795	7.78 ng/ul	80966100	1,4-Dichlorobenzene-d4	7.925

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



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Instrument :
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ClientSampleId :
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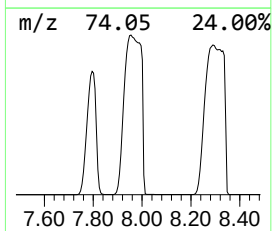
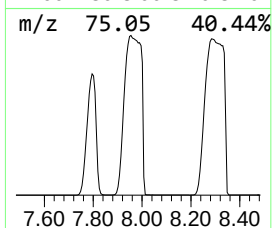
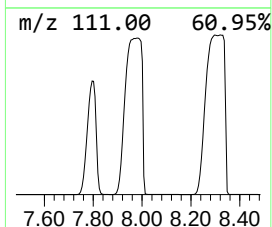
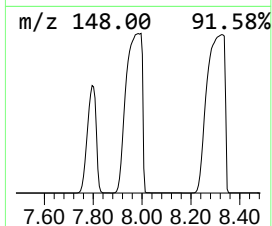
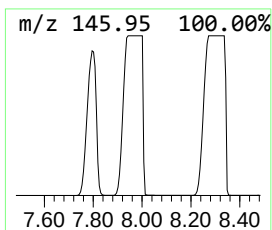
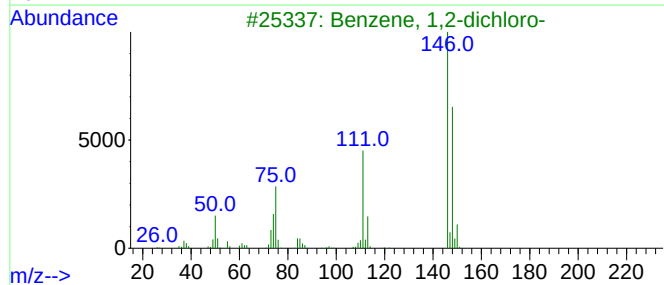
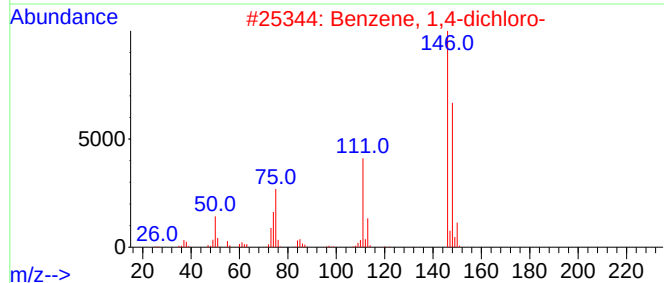
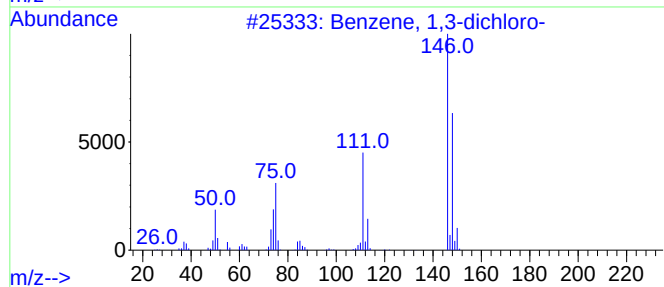
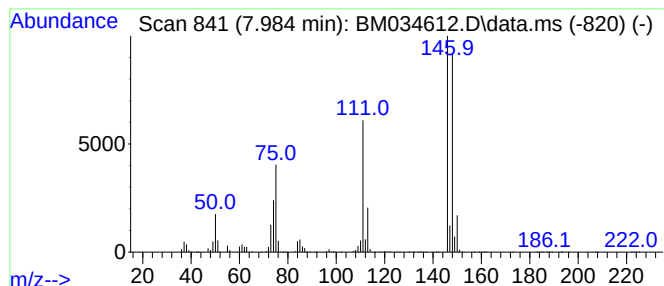
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.984	20.00 ng/ul	208171000	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
4			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95



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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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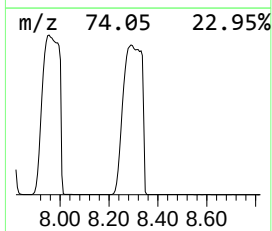
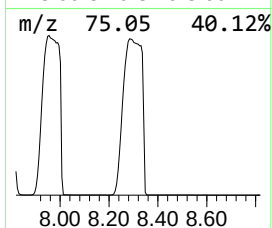
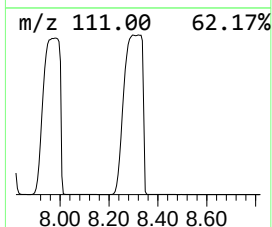
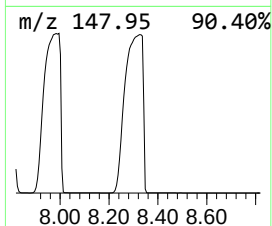
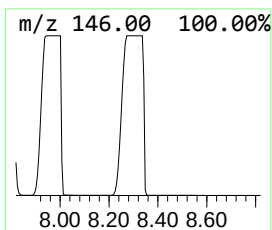
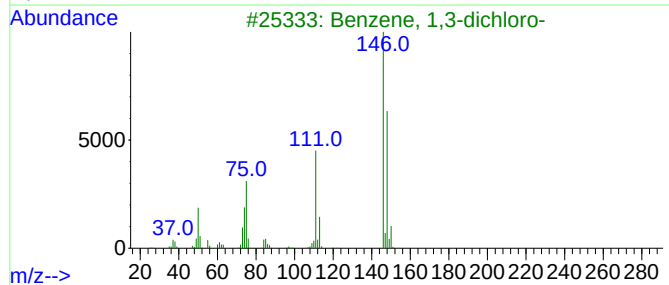
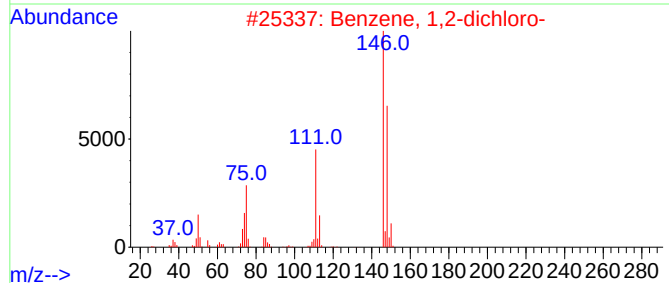
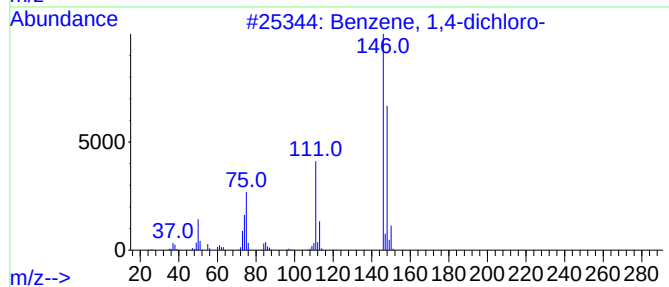
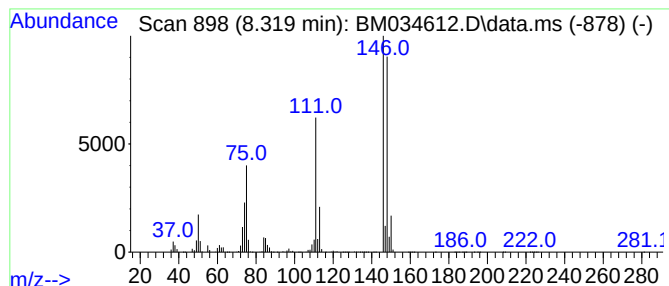
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.319	22.18 ng/ul	230838000	1,4-Dichlorobenzene-d4	7.925

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
3			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95



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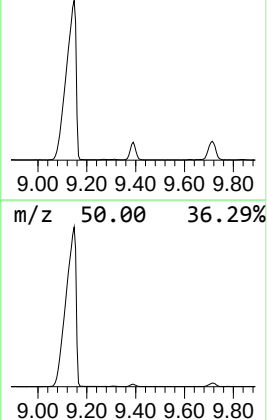
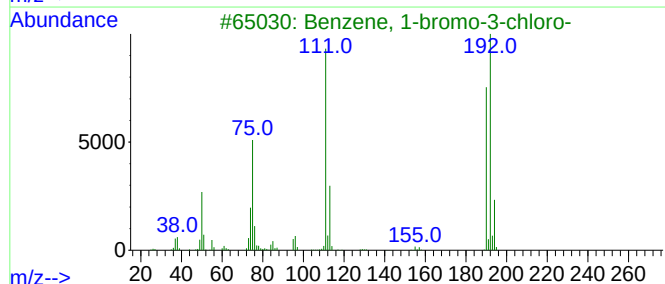
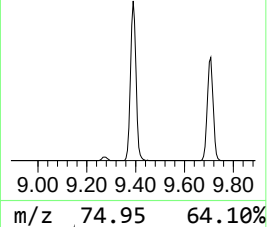
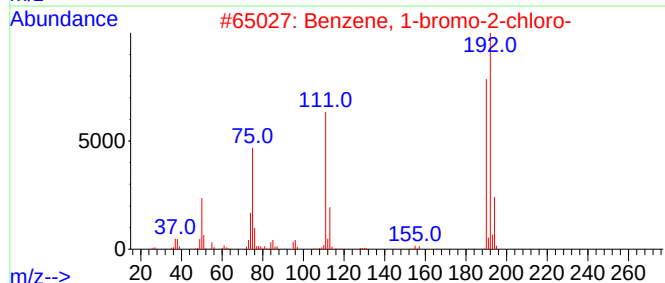
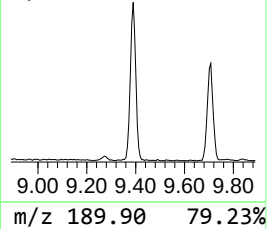
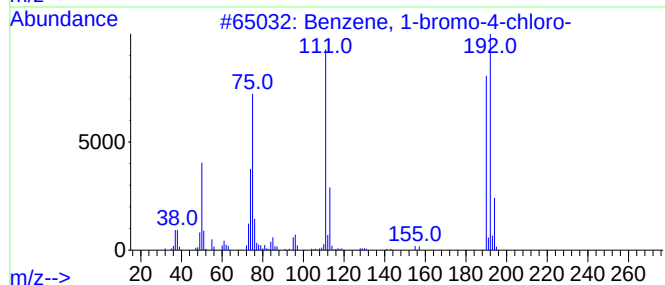
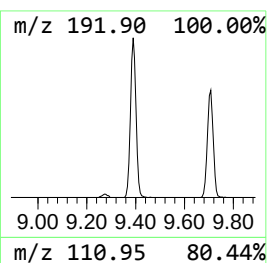
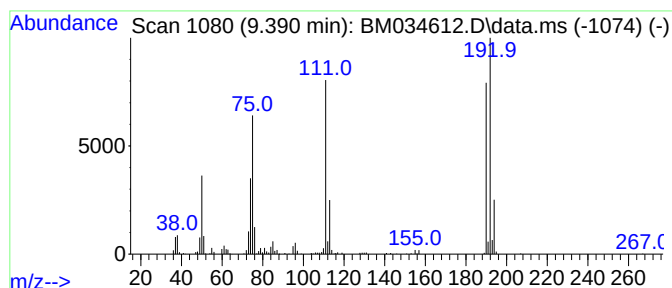
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.390	13.23 ng/ul	658500	Naphthalene-d8	10.707

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	99
2			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4			4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
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Acq On : 11 Apr 2022 16:27
Operator : CG/JU
Sample : N2338-02
Misc :
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Instrument :
BNA_M
ClientSampleId :
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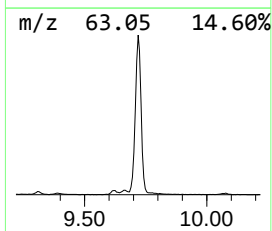
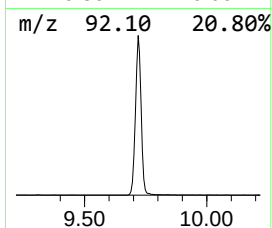
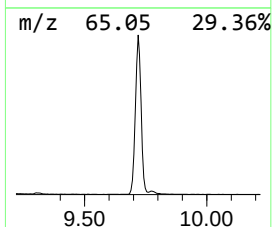
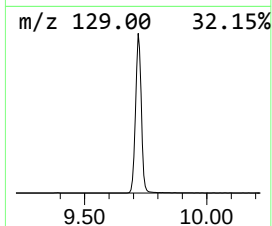
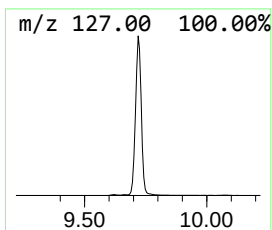
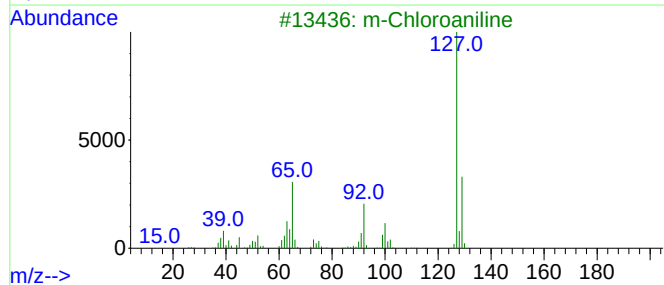
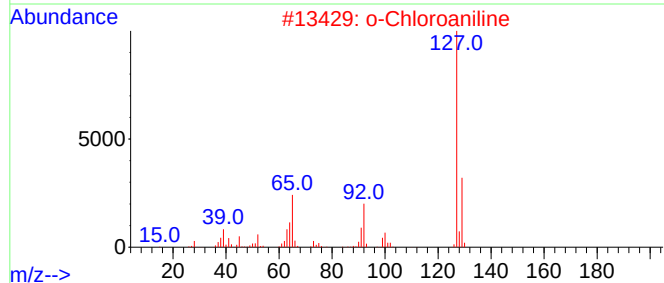
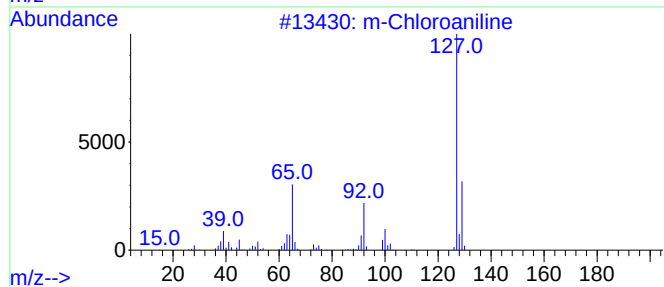
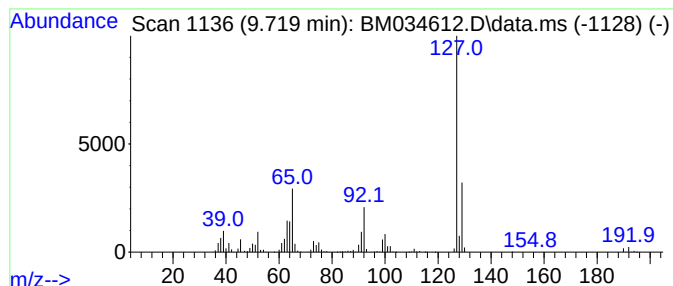
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 m-Chloroaniline Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.719	110.33 ng/ul	5492990	Naphthalene-d8	10.707

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			m-Chloroaniline	127	C6H6ClN	000108-42-9	97
2			o-Chloroaniline	127	C6H6ClN	000095-51-2	96
3			m-Chloroaniline	127	C6H6ClN	000108-42-9	96
4			o-Chloroaniline	127	C6H6ClN	000095-51-2	95
5			o-Chloroaniline	127	C6H6ClN	000095-51-2	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
Operator : CG/JU
Sample : N2338-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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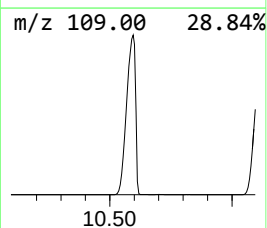
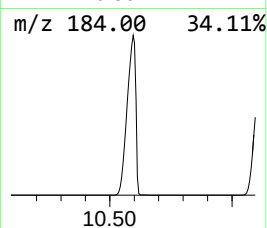
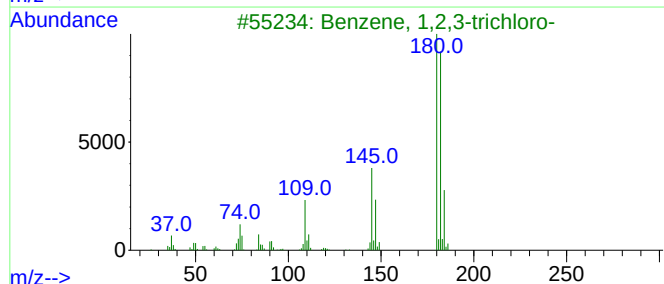
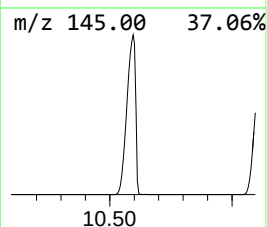
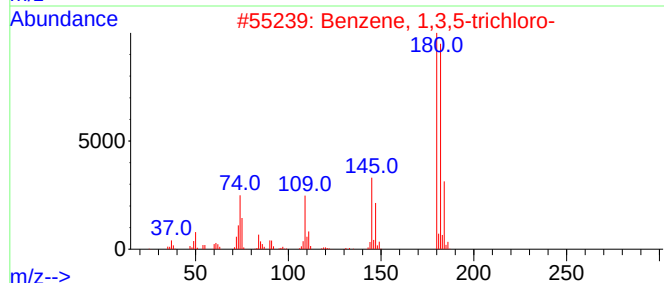
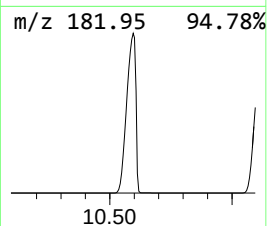
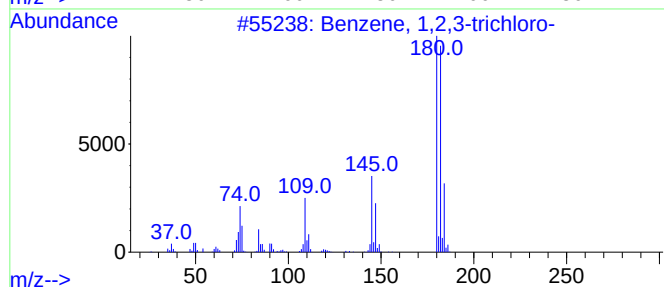
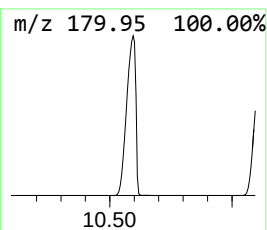
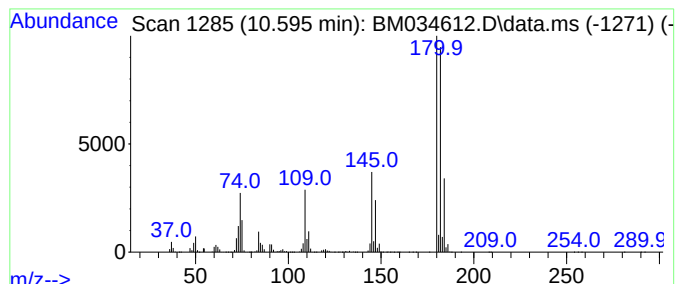
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,2,3-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.595	2101.66 ng/ul	104634000	Naphthalene-d8	10.707

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
Operator : CG/JU
Sample : N2338-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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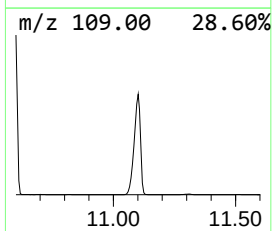
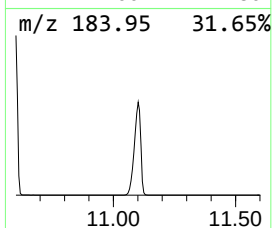
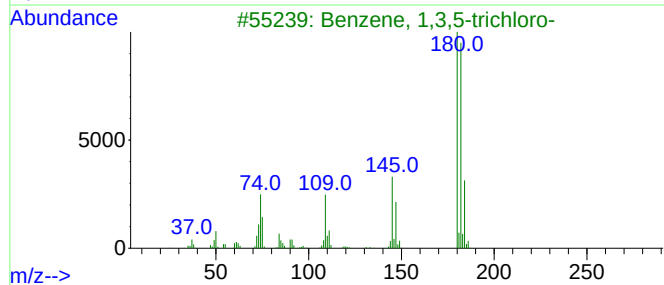
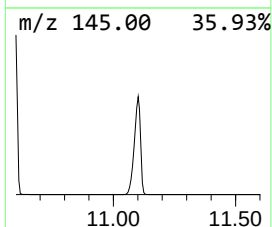
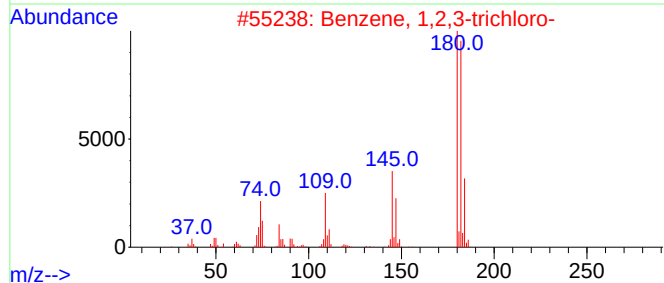
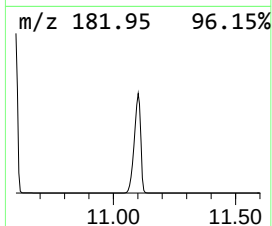
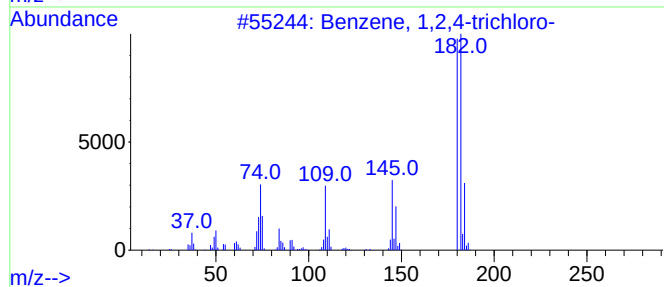
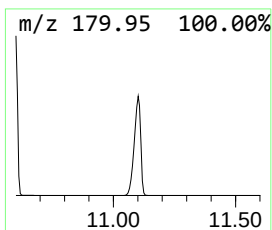
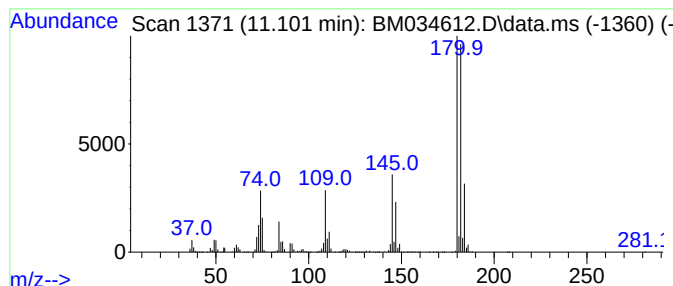
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,4-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.101	906.80 ng/ul	45146200	Naphthalene-d8	10.707

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
Operator : CG/JU
Sample : N2338-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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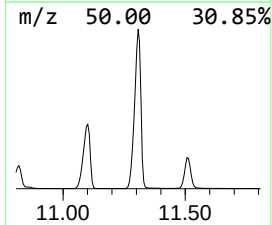
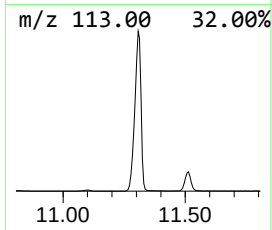
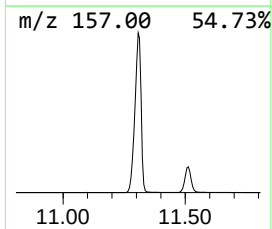
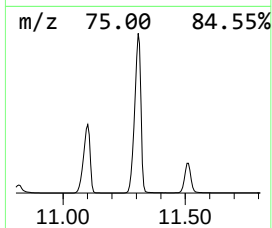
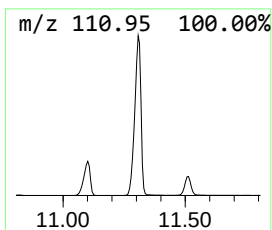
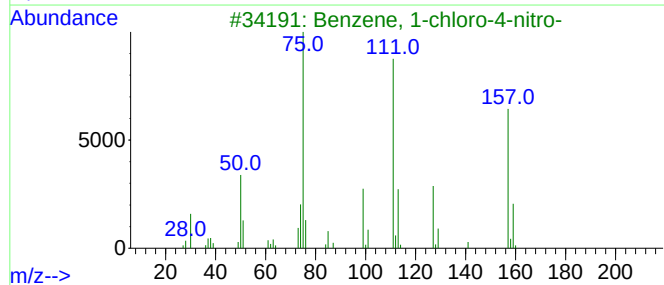
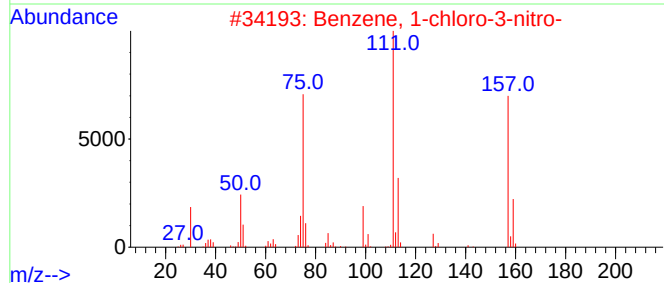
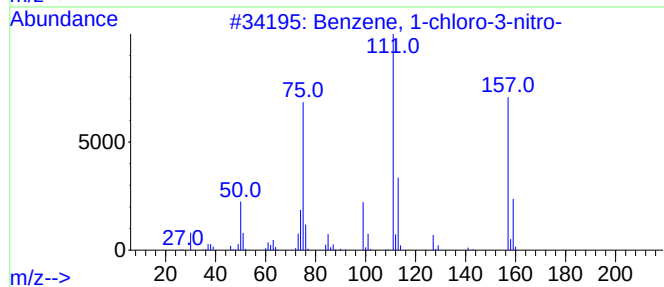
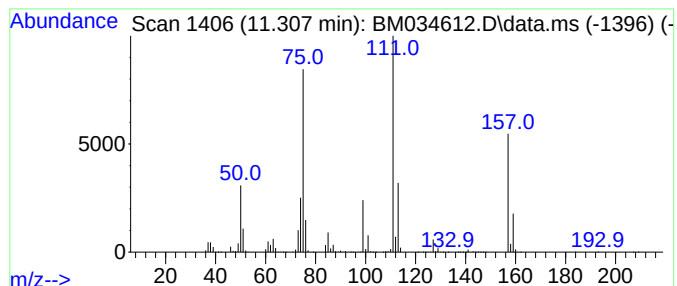
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.307	368.79 ng/ul	18360400	Naphthalene-d8	10.707

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
3			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
Operator : CG/JU
Sample : N2338-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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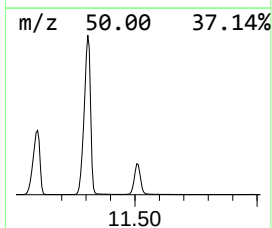
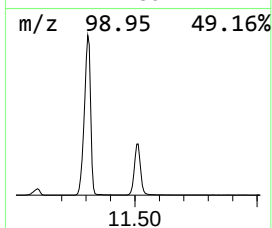
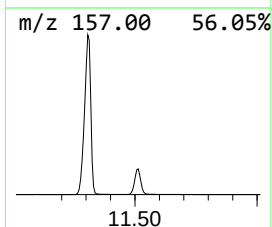
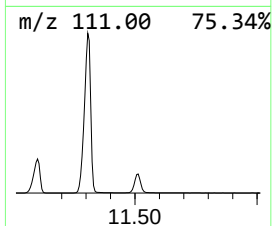
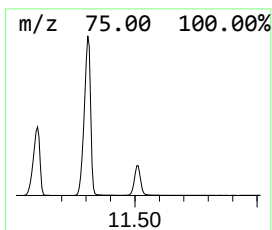
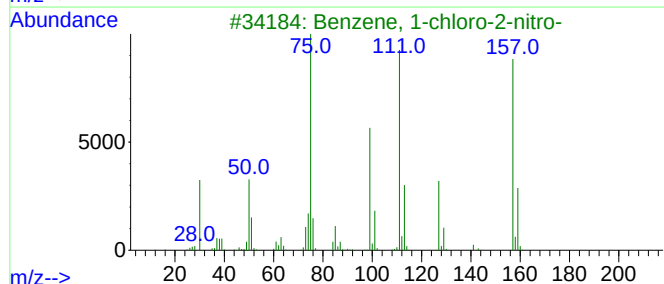
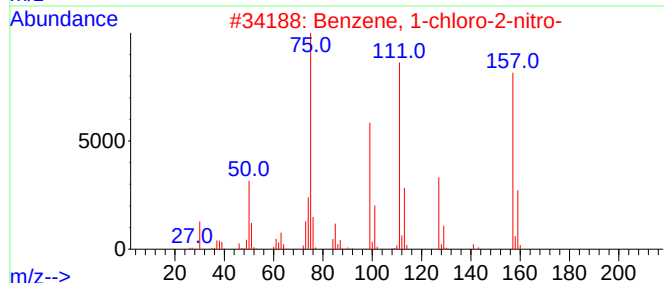
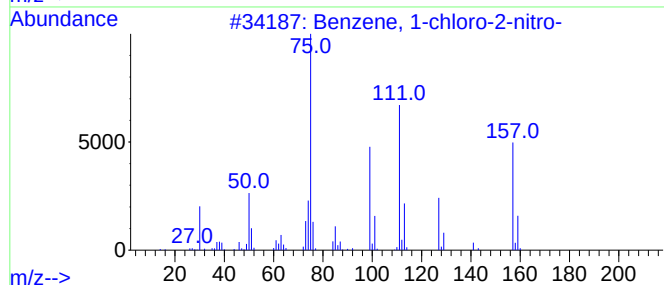
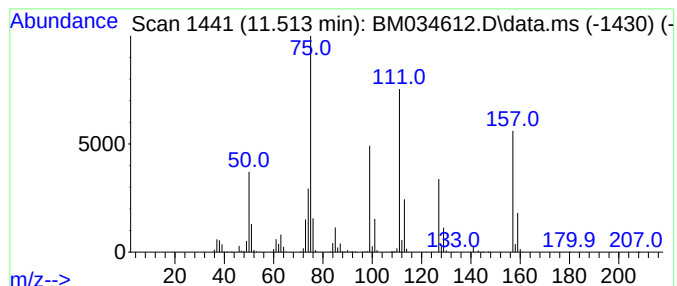
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.513	68.86 ng/ul	3428070	Naphthalene-d8	10.707

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
2			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
3			Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	96
4			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
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Misc :
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Instrument :
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ClientSampleId :
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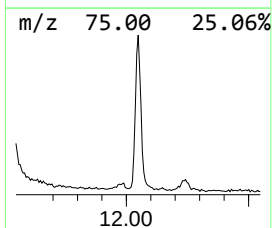
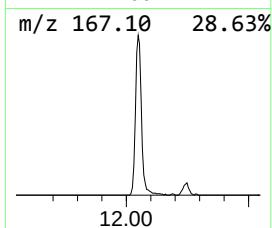
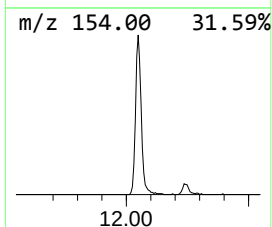
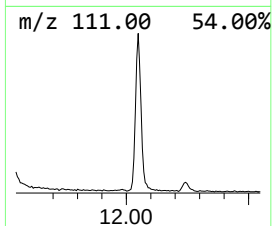
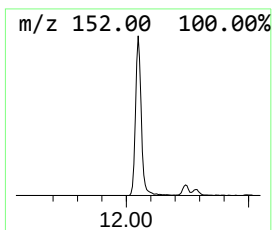
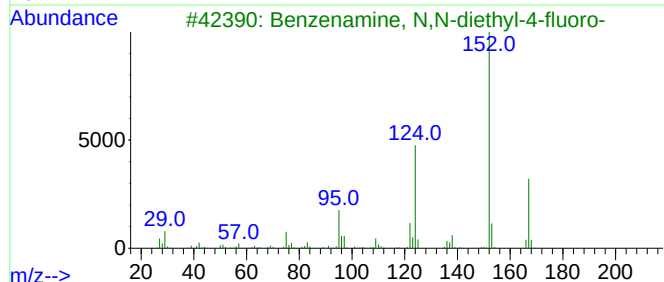
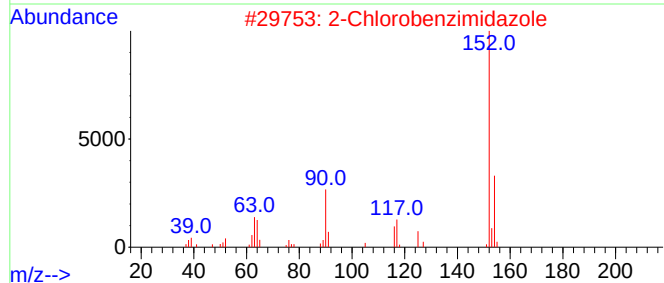
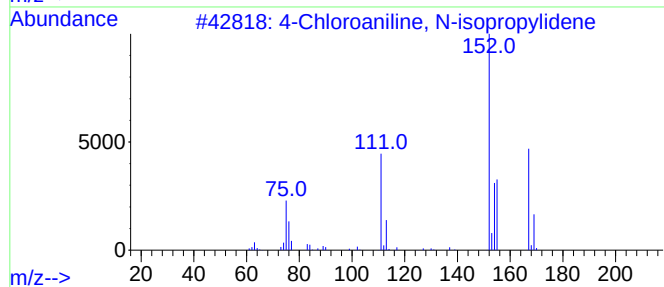
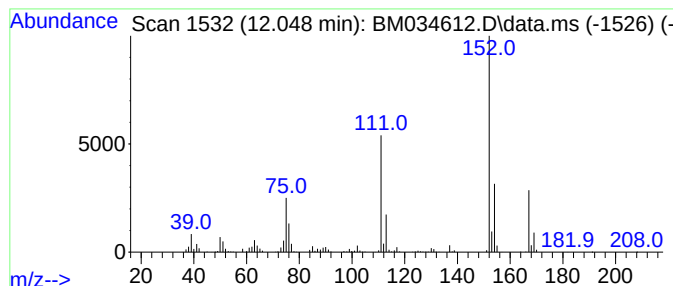
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 4-Chloroaniline, N-isopropy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.048	12.96 ng/ul	645241	Naphthalene-d8	10.707

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloroaniline, N-isopropylidene	167	C9H10ClN	040938-43-0	94
2			2-Chlorobenzimidazole	152	C7H5ClN2	004857-06-1	45
3			Benzenamine, N,N-diethyl-4-fluoro-	167	C10H14FN	000347-39-7	38
4			Acenaphthylene	152	C12H8	000208-96-8	38
5			Benzonitrile, 2-amino-5-chloro-	152	C7H5ClN2	005922-60-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
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Sample : N2338-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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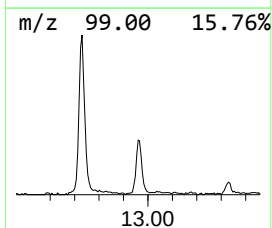
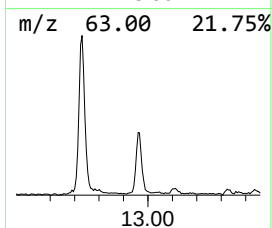
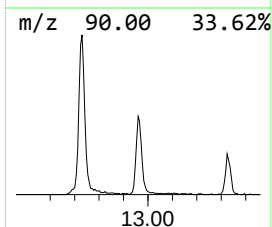
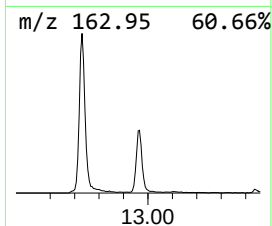
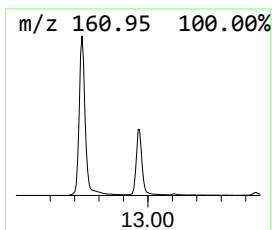
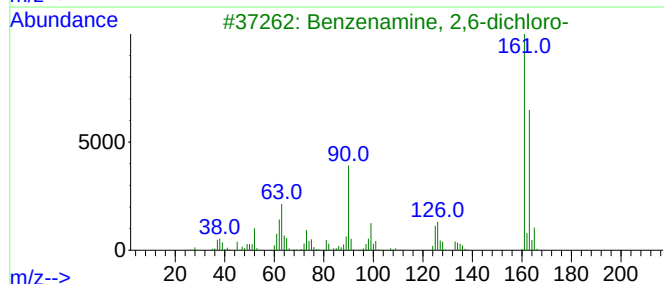
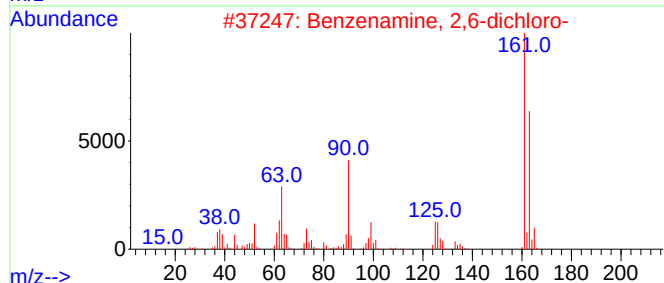
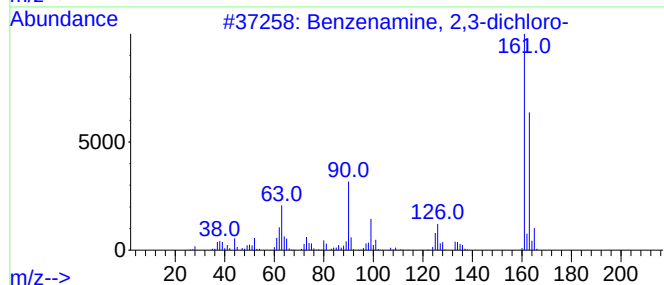
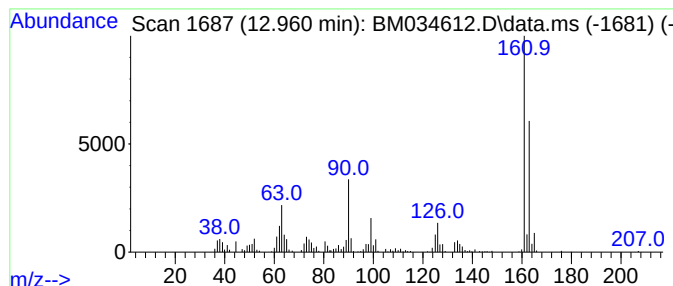
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzenamine, 2,3-dichloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.960	6.45 ng/ul	409084	Acenaphthene-d10	14.525

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	98
2			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	97
3			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	97
4			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97
5			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041122\
Data File : BM034612.D
Acq On : 11 Apr 2022 16:27
Operator : CG/JU
Sample : N2338-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0J71

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.196	13.2	ng/ul	137457000	1	7.925	208171000	20.0
Benzene, 1,4-di...	7.795	7.8	ng/ul	80966100	1	7.925	208171000	20.0
Benzene, 1,3-di...	7.984	20.0	ng/ul	208171000	1	7.925	208171000	20.0
Benzene, 1,2-di...	8.319	22.2	ng/ul	230838000	1	7.925	208171000	20.0
Benzene, 1-brom...	9.390	13.2	ng/ul	658500	2	10.707	995725	20.0
m-Chloroaniline	9.719	110.3	ng/ul	5492990	2	10.707	995725	20.0
Benzene, 1,2,3-...	10.595	2101.7	ng/ul	104634000	2	10.707	995725	20.0
Benzene, 1,2,4-...	11.101	906.8	ng/ul	45146200	2	10.707	995725	20.0
Benzene, 1-chlo...	11.307	368.8	ng/ul	18360400	2	10.707	995725	20.0
Benzene, 1-chlo...	11.513	68.9	ng/ul	3428070	2	10.707	995725	20.0
4-Chloroaniline...	12.048	13.0	ng/ul	645241	2	10.707	995725	20.0
Benzenamine, 2,...	12.960	6.5	ng/ul	409084	3	14.525	1269240	20.0