

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
 Data File : BP020094.D
 Acq On : 23 Apr 2024 04:20
 Operator : MA/JU
 Sample : P2184-02
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CORE5

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: 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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020094.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.652	92	96	123	rVB	107324	347228	0.13%	0.039%
2	5.093	320	341	343	rBV3	34637891	192115805	73.19%	21.358%
3	5.246	366	367	369	rVB	33908574	12704599	4.84%	1.412%
4	6.358	550	556	566	rBV	484683	826102	0.31%	0.092%
5	6.734	614	620	632	rVB	309993	649229	0.25%	0.072%
6	6.840	632	638	653	rVB2	202967	555713	0.21%	0.062%
7	7.040	655	672	673	rBV2	1631040	4874925	1.86%	0.542%
8	7.228	692	704	723	rVB3	401352	1942810	0.74%	0.216%
9	7.558	750	760	761	rBV	8403424	16500687	6.29%	1.834%
10	7.822	778	805	806	rBV5	25714343	201773587	76.87%	22.432%
11	8.181	837	866	867	rBV4	33958332	262496239	100.00%	29.183%
12	8.381	894	900	915	rVB	644538	1128274	0.43%	0.125%
13	8.834	970	977	993	rBV	1154289	1956880	0.75%	0.218%
14	9.152	1024	1031	1041	rVB	699667	1134454	0.43%	0.126%
15	9.475	1077	1086	1092	rVV	4031922	7356126	2.80%	0.818%
16	9.534	1092	1096	1109	rVB	961777	1561320	0.59%	0.174%
17	10.069	1179	1187	1193	rBV	1168045	2150446	0.82%	0.239%
18	10.134	1193	1198	1206	rVB	496008	872770	0.33%	0.097%
19	10.228	1206	1214	1220	rBV	1618832	3461065	1.32%	0.385%
20	10.340	1220	1233	1237	rVV2	30133561	83328217	31.74%	9.264%
21	10.446	1245	1251	1260	rVB	961765	1461899	0.56%	0.163%
22	10.599	1262	1277	1292	rBV	14785399	43197096	16.46%	4.802%
23	10.828	1306	1316	1326	rVB	12316913	22300889	8.50%	2.479%
24	12.498	1585	1600	1610	rVV2	1043526	2215889	0.84%	0.246%
25	12.728	1628	1639	1648	rBV	533687	887966	0.34%	0.099%
26	13.087	1695	1700	1705	rVV	462218	737508	0.28%	0.082%
27	13.340	1736	1743	1749	rBV	493076	908946	0.35%	0.101%
28	13.663	1789	1798	1803	rBV	194701	339862	0.13%	0.038%
29	13.734	1803	1810	1825	rVB	1824252	3052752	1.16%	0.339%
30	14.004	1850	1856	1871	rVB	2189516	3457142	1.32%	0.384%
31	14.322	1903	1910	1918	rBV	1163033	1725034	0.66%	0.192%
32	15.345	2075	2084	2098	rVB	2399567	3841834	1.46%	0.427%
33	15.498	2103	2110	2120	rBV	453840	730360	0.28%	0.081%
34	17.175	2388	2395	2403	rBV	1119108	1669716	0.64%	0.186%
35	17.269	2405	2411	2423	rBV	2595560	3649557	1.39%	0.406%
36	19.675	2813	2820	2826	rBV	2665779	4186422	1.59%	0.465%

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ClientSampleId :
C0RE5

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-BP041924.MA.M
Title : SVOA CALIBRATION

37	21.663	3153	3158	3169	rVB2	967717	1702108	0.65%	0.189%
38	24.821	3686	3695	3708	rVB2	1478537	3987707	1.52%	0.443%
39	25.057	3726	3735	3750	rVB2	517160	1699340	0.65%	0.189%

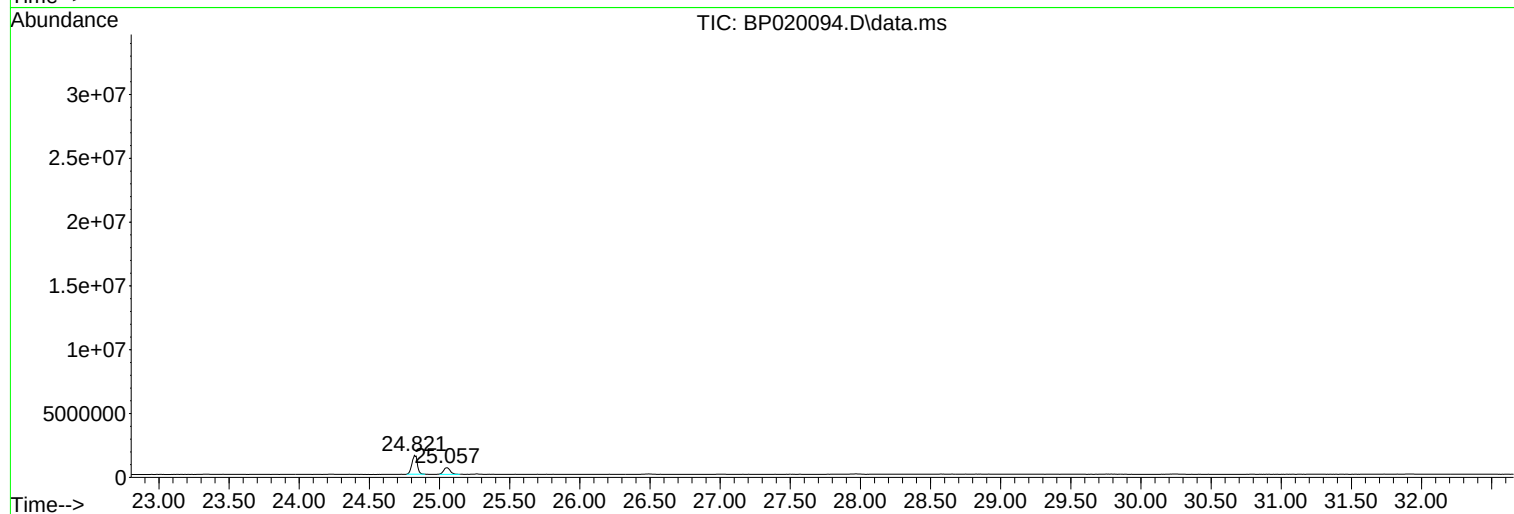
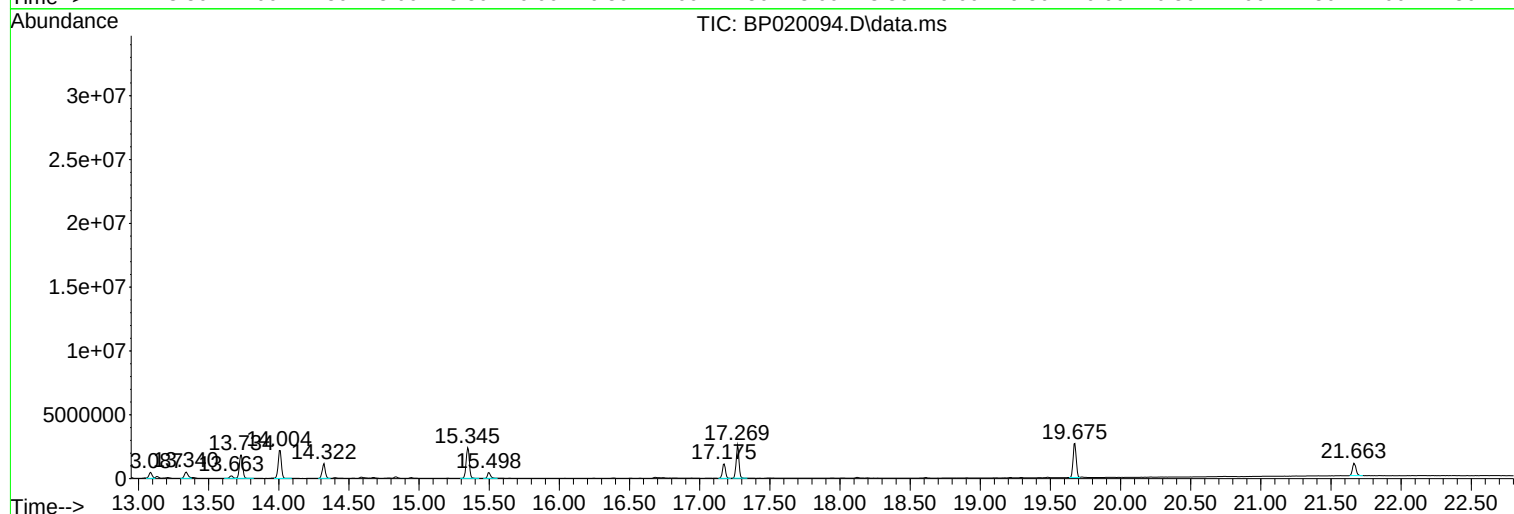
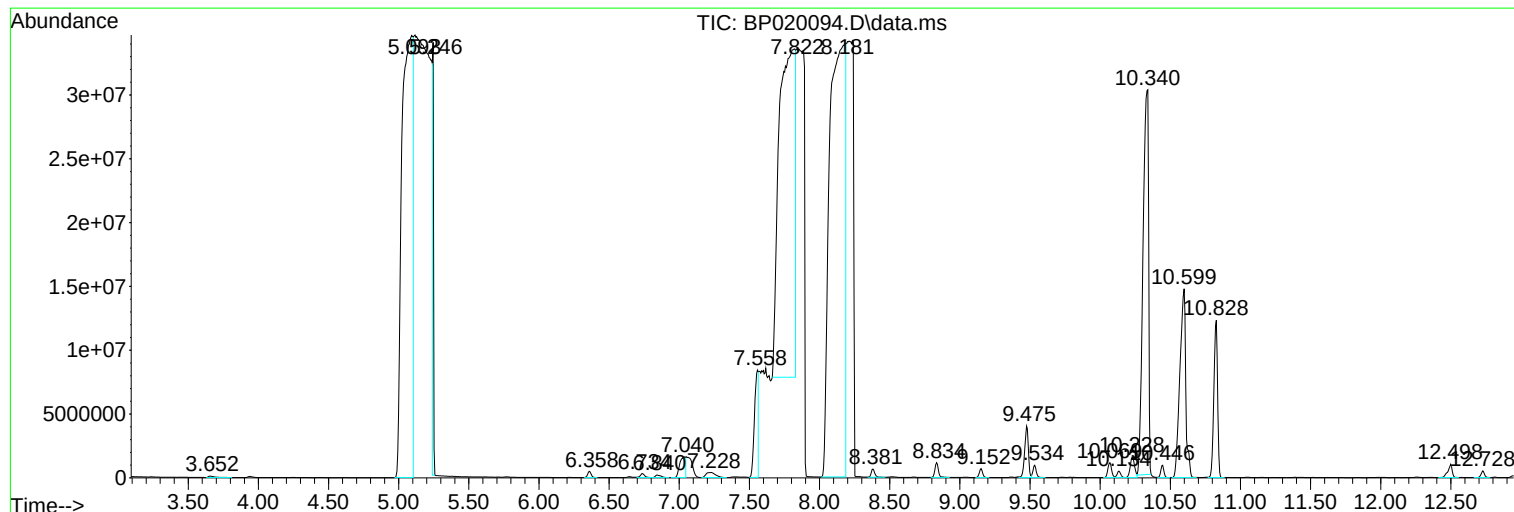
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ClientSampleId :
CORE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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



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Operator : MA/JU
Sample : P2184-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CORE5

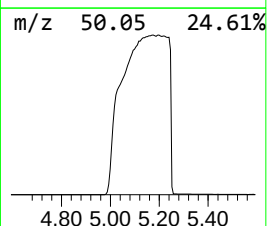
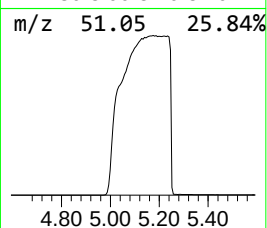
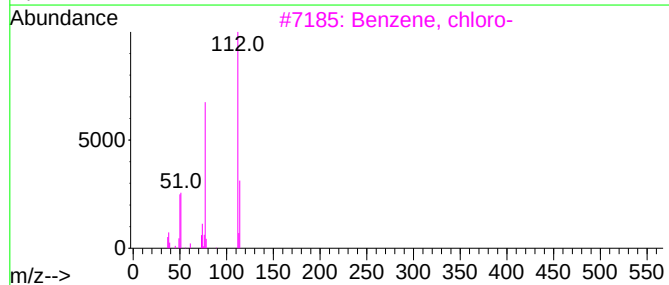
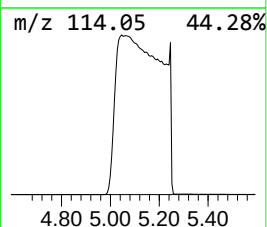
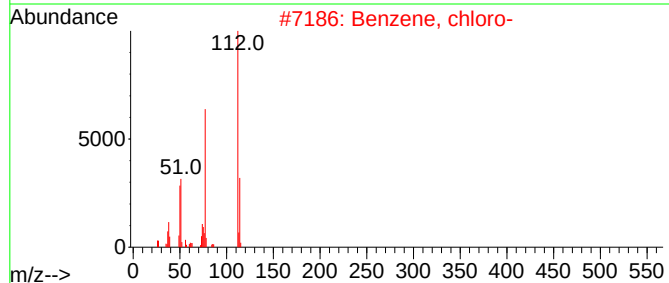
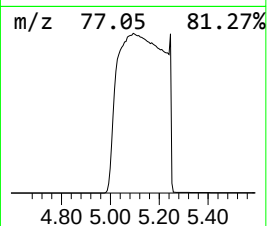
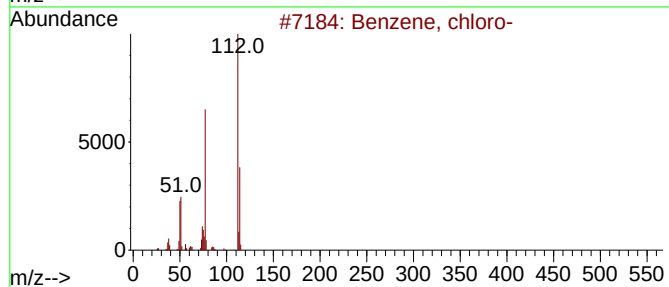
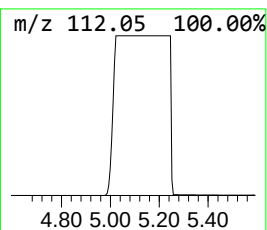
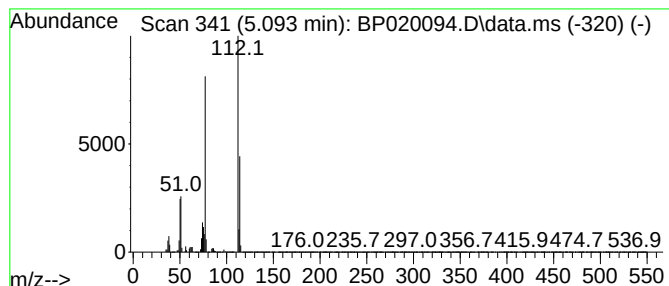
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.093	19.04 ng/ul	192116000	1,4-Dichlorobenzene-d4	7.705

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



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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CORE5

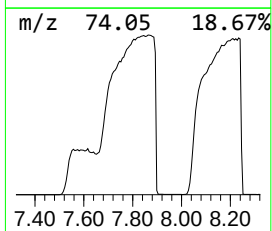
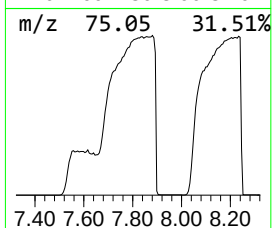
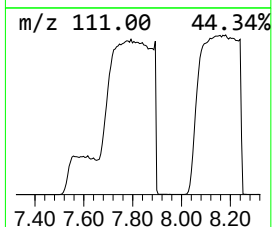
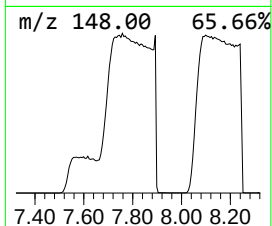
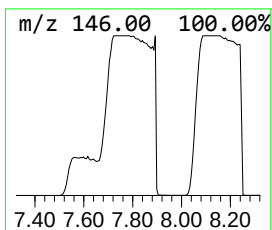
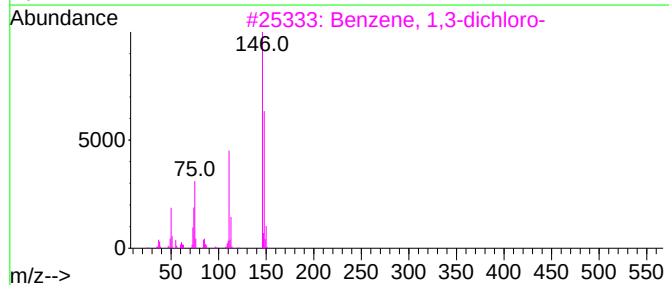
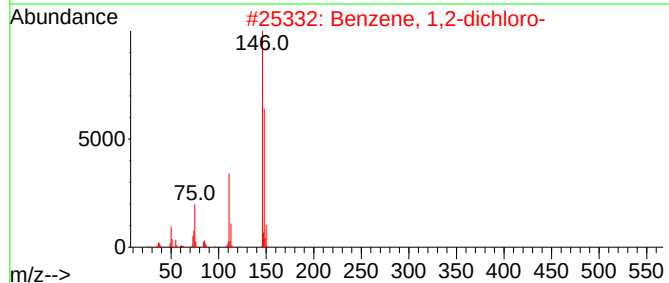
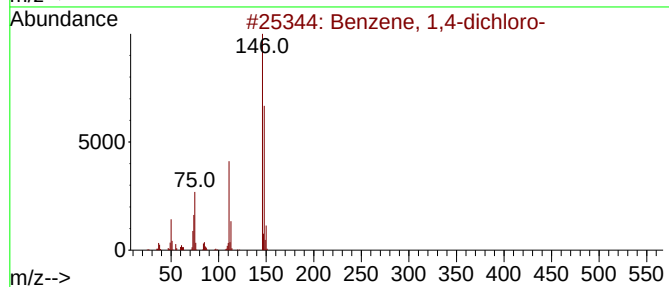
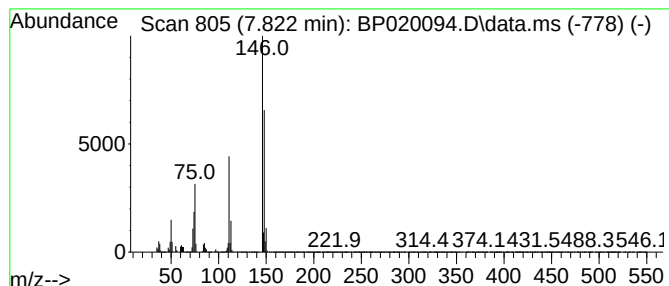
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.822	20.00 ng/ul	201774000	1,4-Dichlorobenzene-d4	7.705

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



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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CORE5

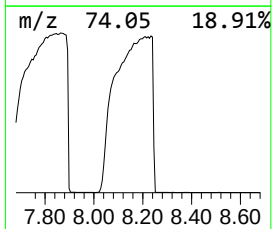
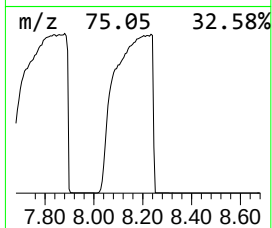
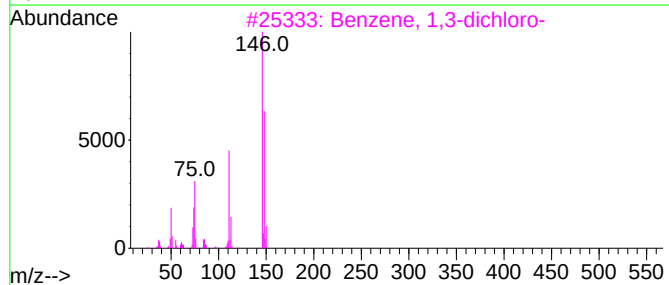
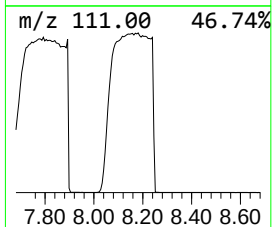
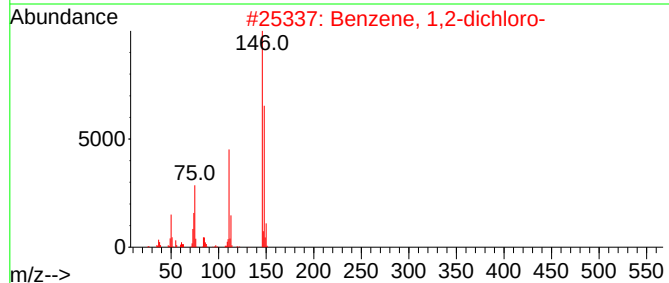
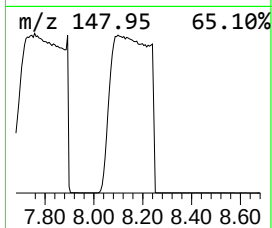
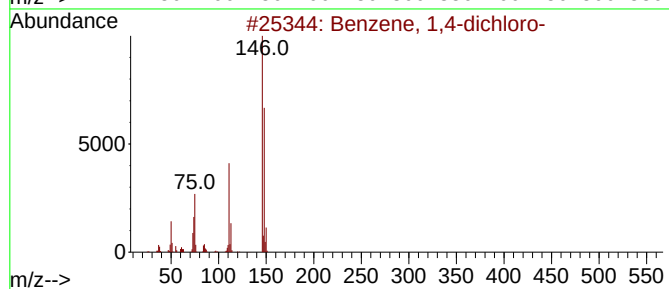
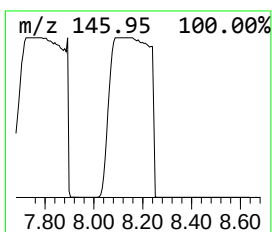
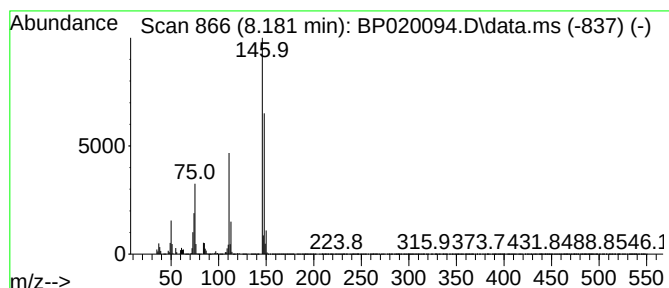
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,2-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.181	26.02 ng/ul	262496000	1,4-Dichlorobenzene-d4	7.705

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



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ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
CORE5

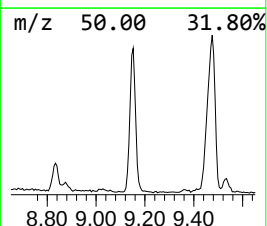
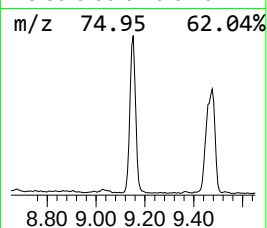
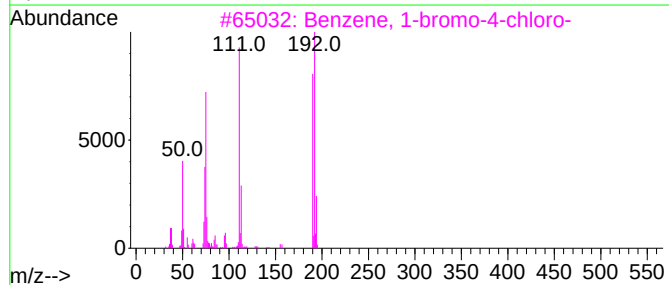
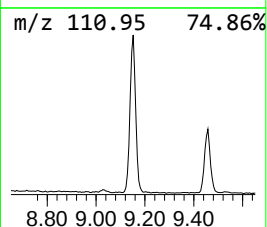
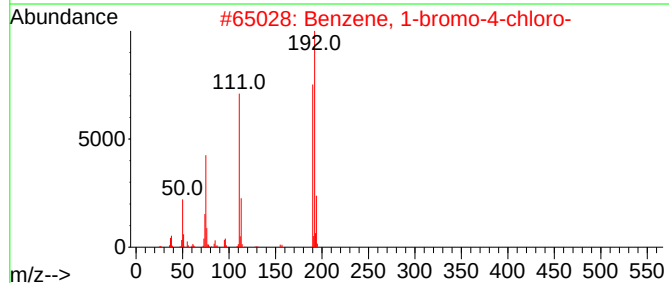
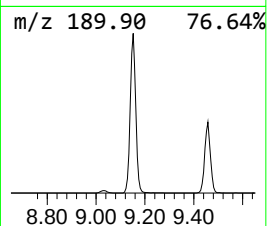
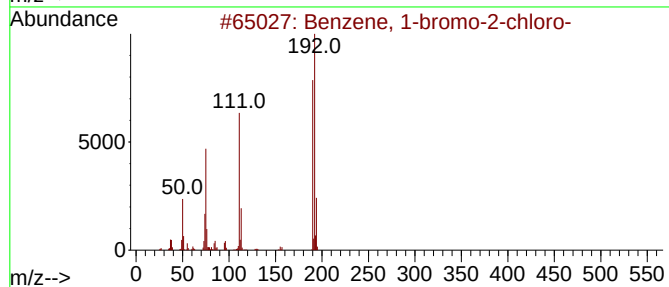
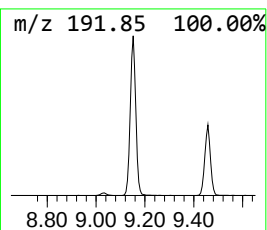
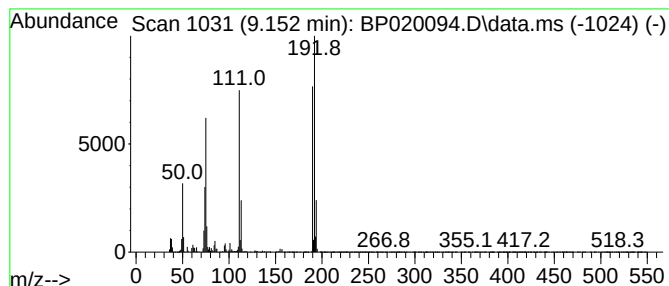
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.152	15.52 ng/ul	1134450	Naphthalene-d8	10.446

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	94



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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CORE5

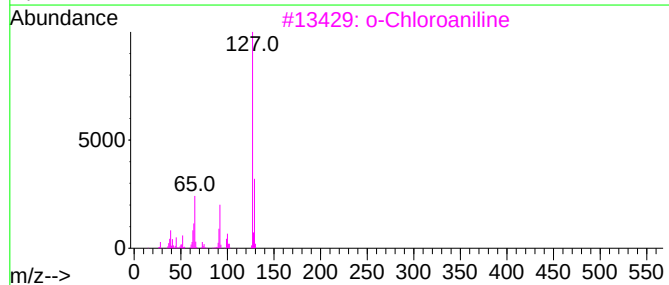
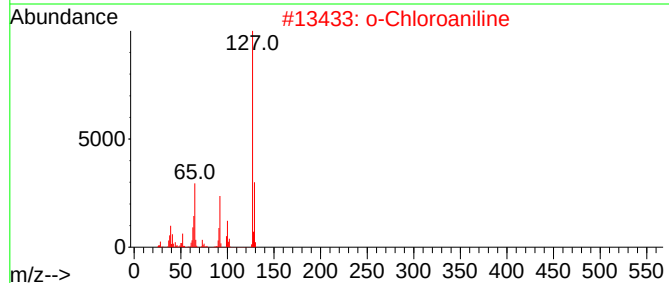
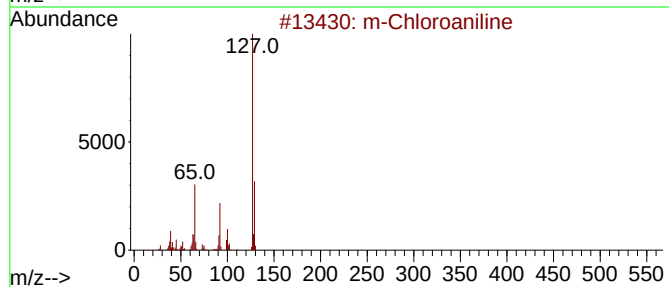
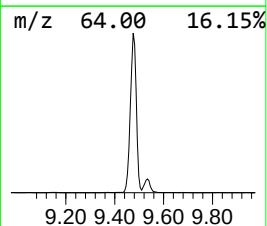
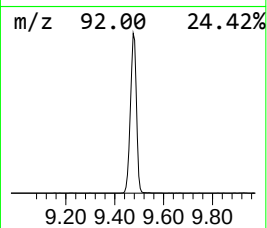
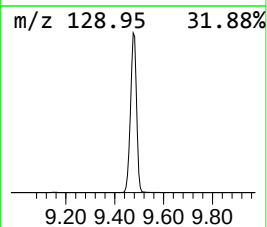
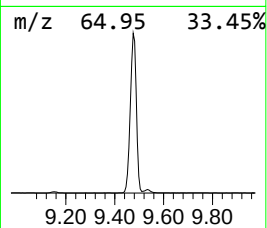
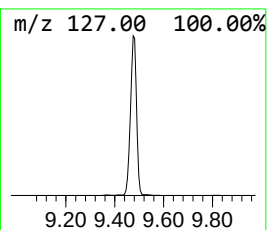
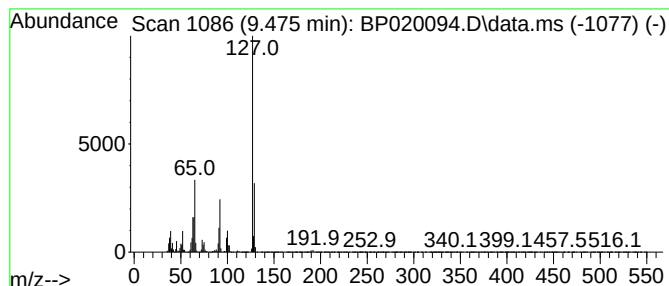
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 m-Chloroaniline Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.475	100.64 ng/ul	7356130	Naphthalene-d8	10.446

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			o-Chloroaniline	127	C6H6ClN	000095-51-2	96
4			o-Chloroaniline	127	C6H6ClN	000095-51-2	95
5			m-Chloroaniline	127	C6H6ClN	000108-42-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020094.D
Acq On : 23 Apr 2024 04:20
Operator : MA/JU
Sample : P2184-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

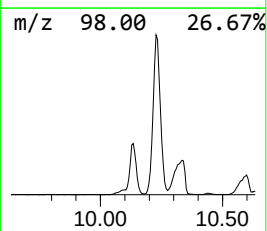
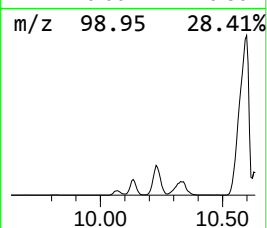
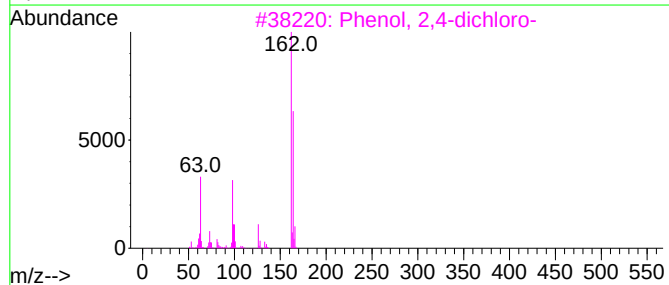
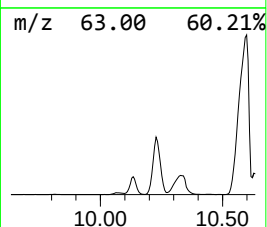
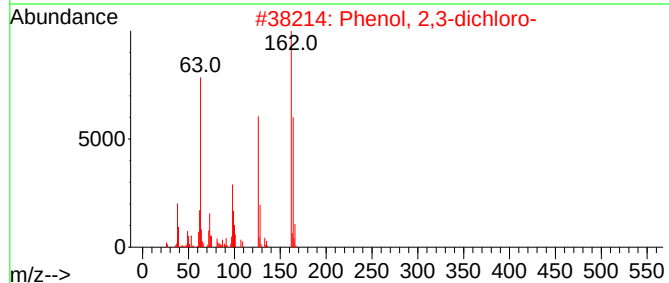
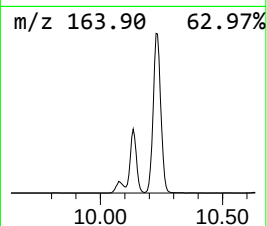
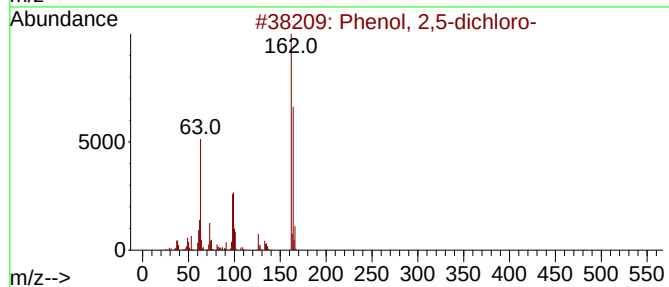
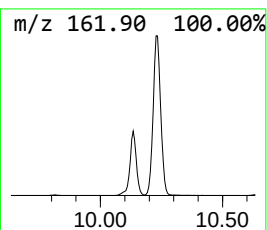
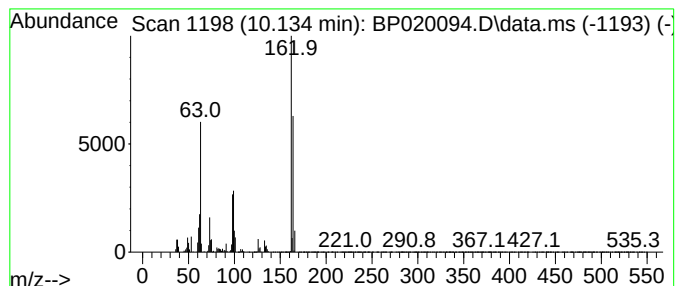
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 2,5-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.134	11.94 ng/ul	872770	Naphthalene-d8	10.446	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,5-dichloro-	162	C6H4Cl2O	000583-78-8	95
2	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	94
3	Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	94
4	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	94
5	Phenol, 2,5-dichloro-	162	C6H4Cl2O	000583-78-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020094.D
Acq On : 23 Apr 2024 04:20
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Sample : P2184-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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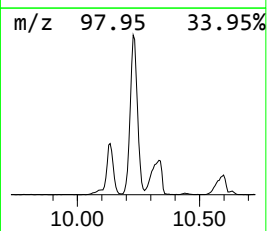
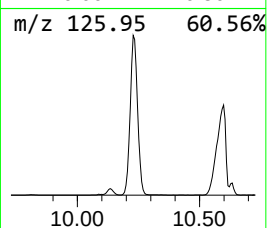
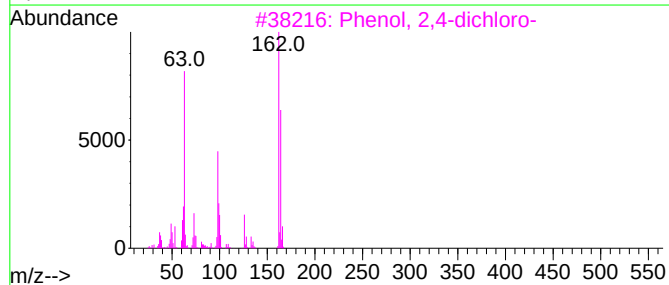
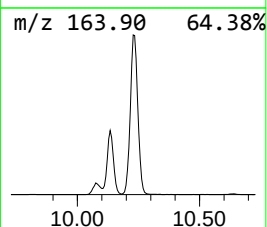
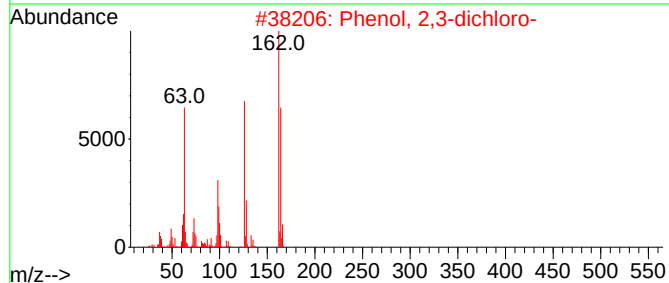
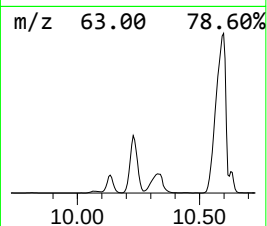
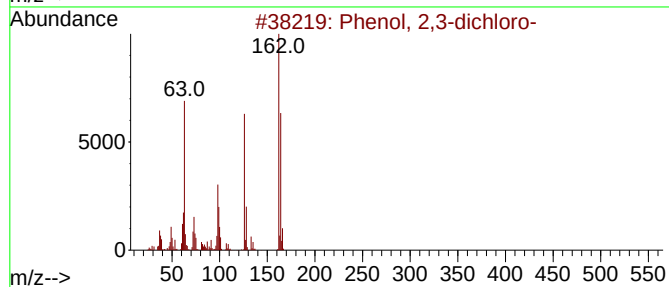
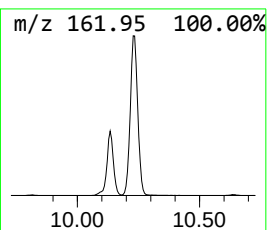
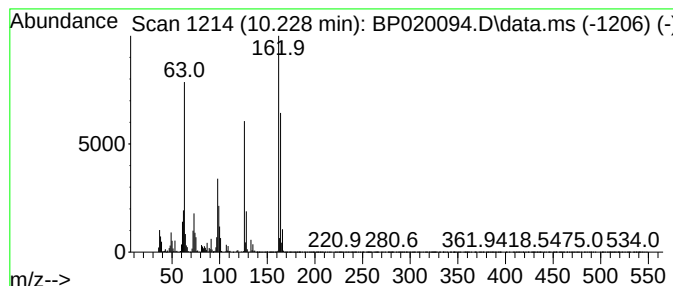
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Phenol, 2,3-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.228	47.35 ng/ul	3461070	Naphthalene-d8	10.446

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,3-dichloro-		162	C6H4Cl2O	000576-24-9	99	
2	Phenol, 2,3-dichloro-		162	C6H4Cl2O	000576-24-9	98	
3	Phenol, 2,4-dichloro-		162	C6H4Cl2O	000120-83-2	98	
4	Phenol, 2,4-dichloro-		162	C6H4Cl2O	000120-83-2	96	
5	Phenol, 2,3-dichloro-		162	C6H4Cl2O	000576-24-9	95	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020094.D
Acq On : 23 Apr 2024 04:20
Operator : MA/JU
Sample : P2184-02
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ClientSampleId :
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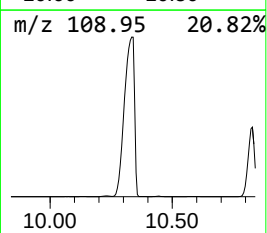
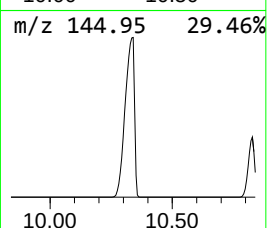
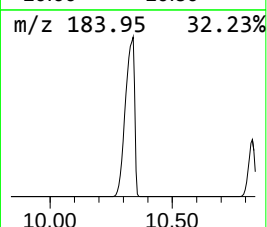
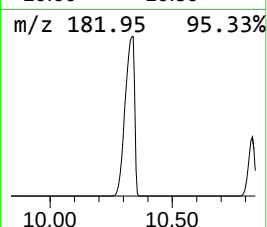
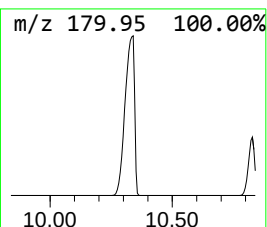
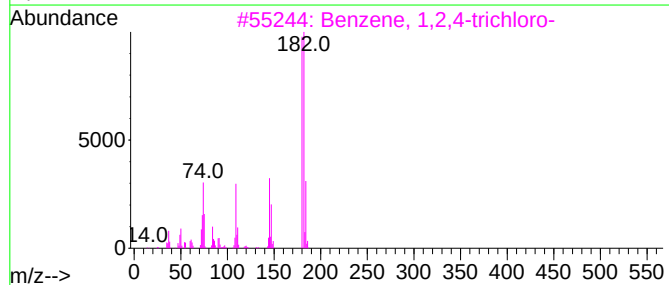
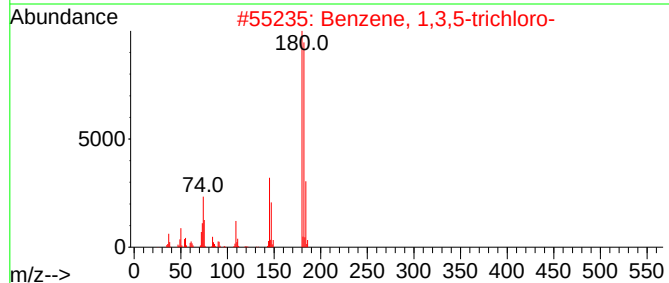
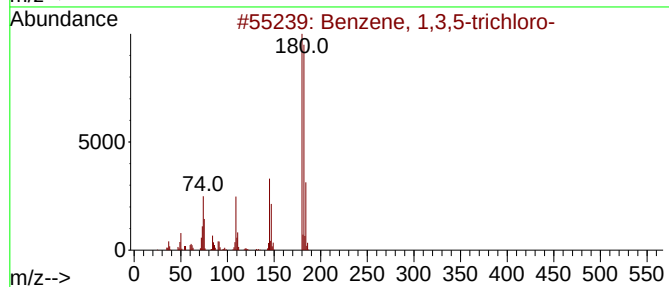
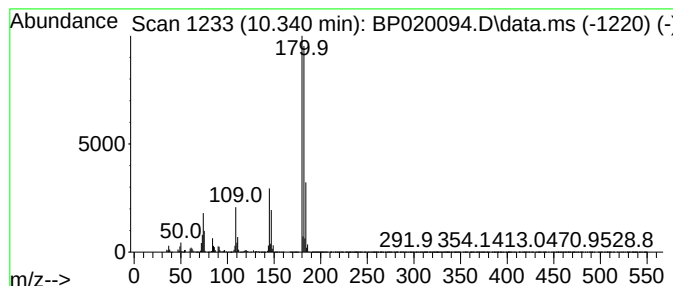
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.340	1140.00 ng/ul	83328200	Naphthalene-d8	10.446

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	97
3			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
4			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
5			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020094.D
Acq On : 23 Apr 2024 04:20
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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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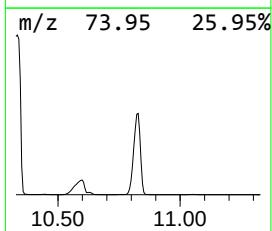
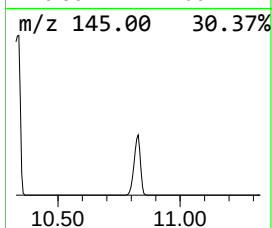
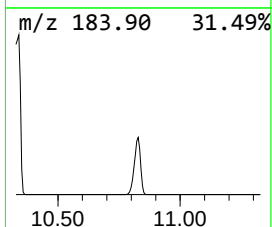
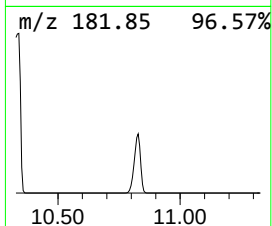
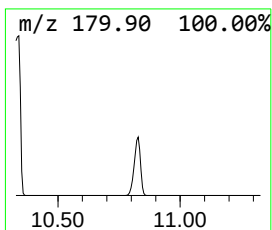
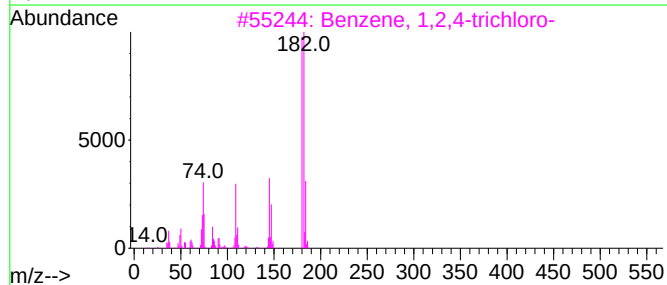
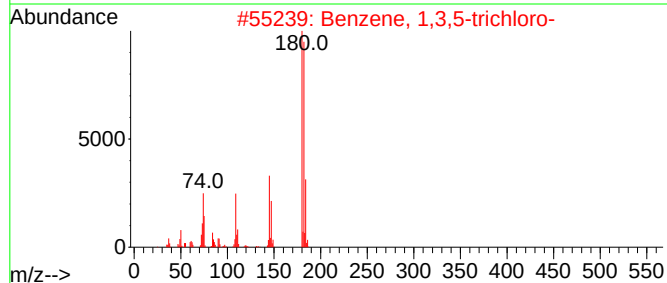
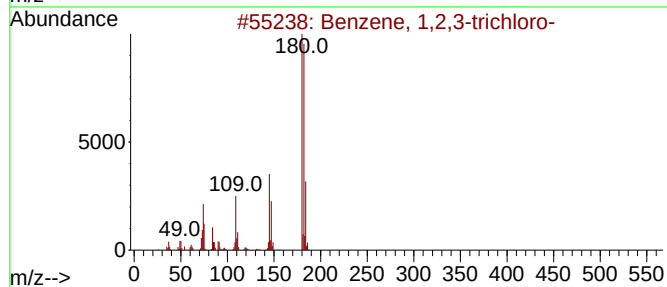
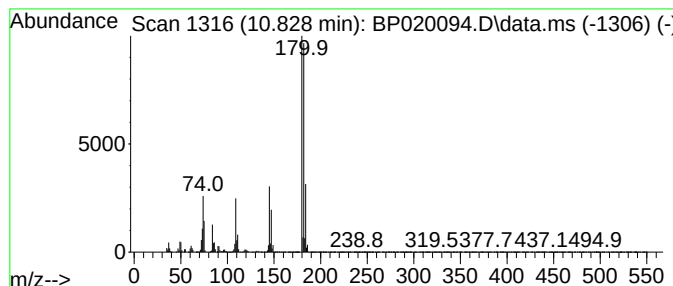
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.828	305.10 ng/ul	22300900	Naphthalene-d8	10.446

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	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_P\Data\BP042224\
Data File : BP020094.D
Acq On : 23 Apr 2024 04:20
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Misc :
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Instrument :
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ClientSampleId :
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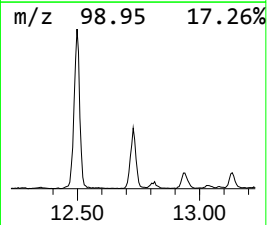
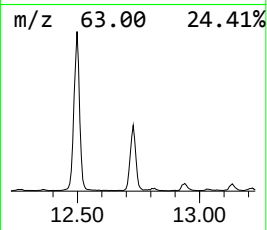
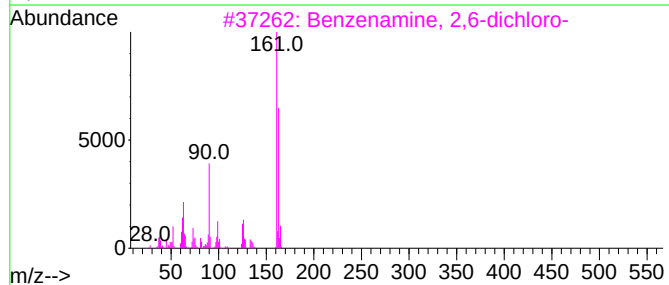
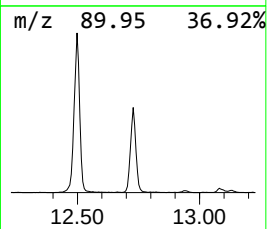
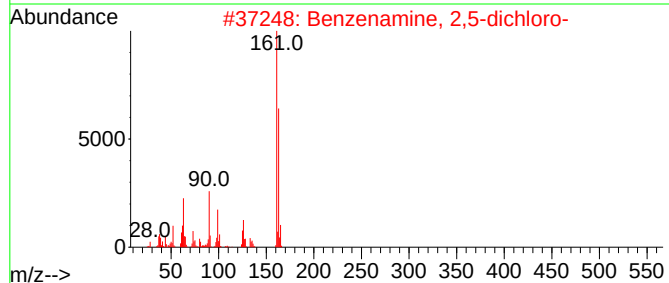
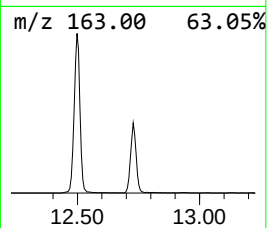
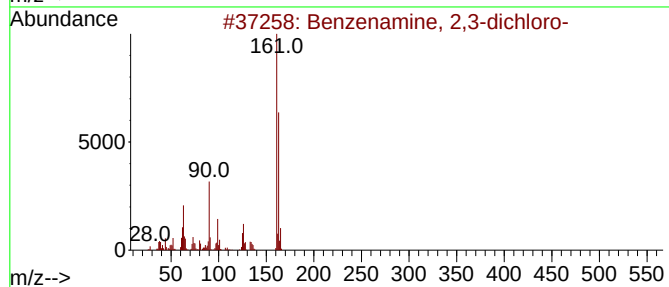
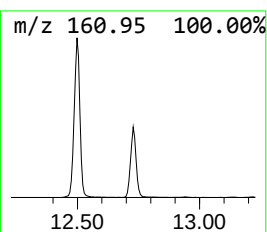
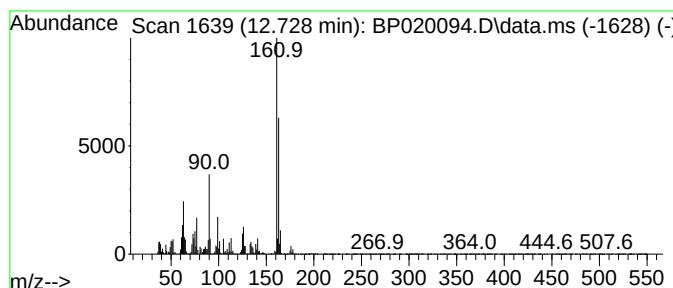
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzenamine, 2,3-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.728	10.30 ng/ul	887966	Acenaphthene-d10	14.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	98
2			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	98
3			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	97
4			Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	97
5			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
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Sample : P2184-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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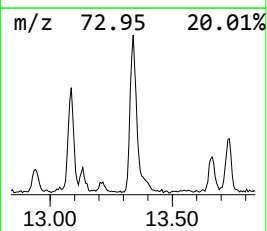
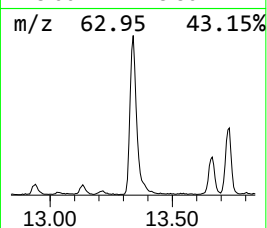
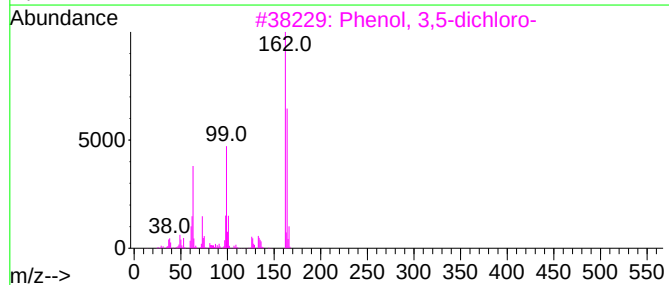
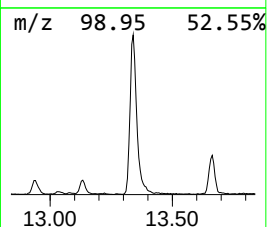
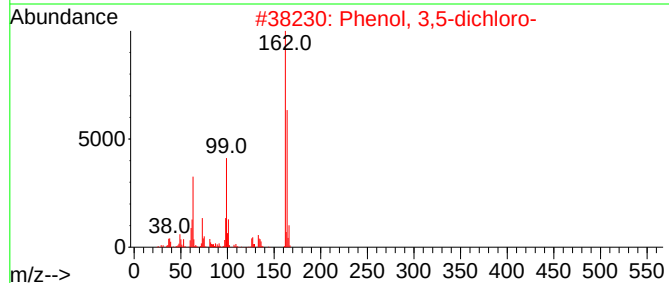
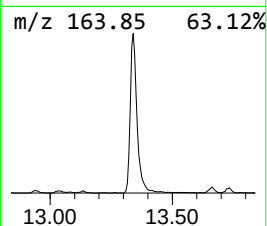
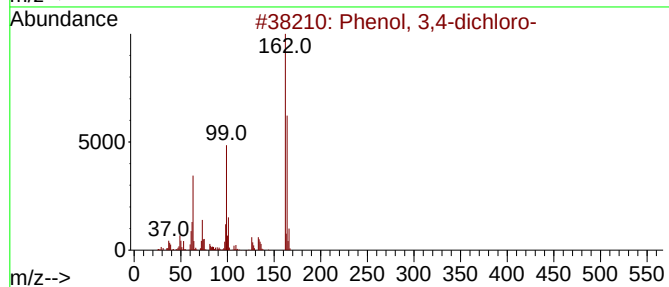
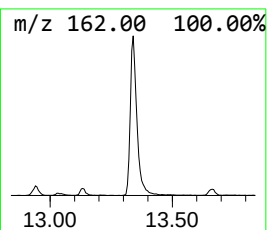
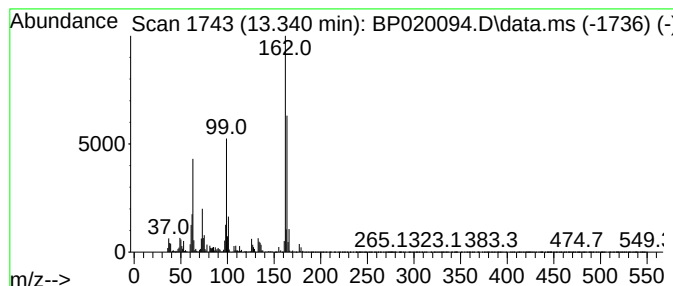
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenol, 3,4-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.340	10.54 ng/ul	908946	Acenaphthene-d10	14.322

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
2			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3			Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
4			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97
5			Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020094.D
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Sample : P2184-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

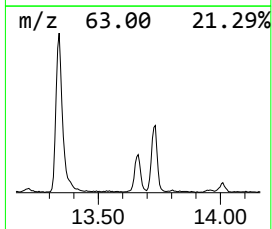
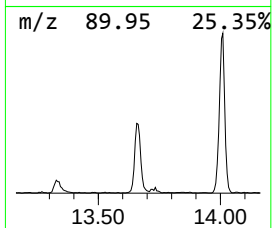
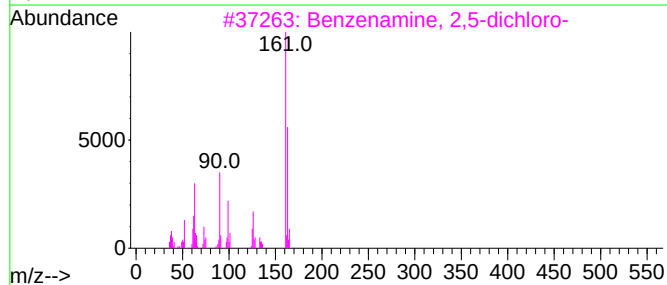
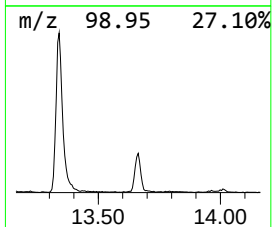
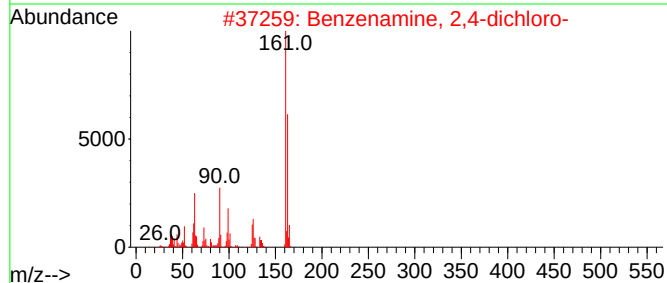
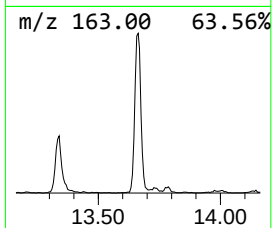
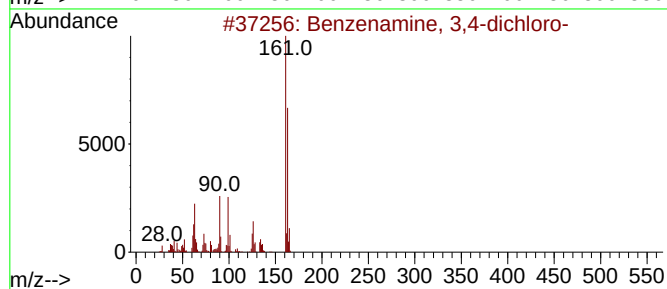
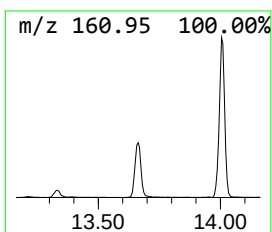
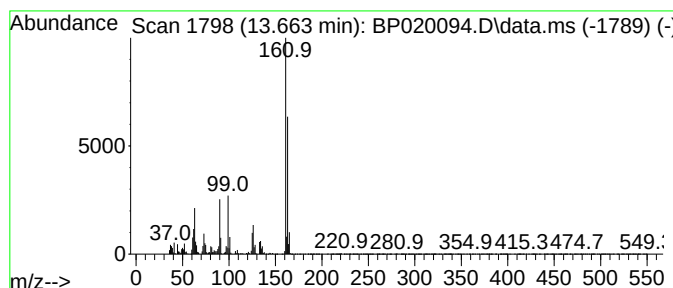
Instrument :
BNA_P
ClientSampleId :
CORE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzenamine, 3,4-dichloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.663	3.94 ng/ul	339862	Acenaphthene-d10	14.322	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3,4-dichloro-	161	C6H5Cl2N	000095-76-1	98
2	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	97
3	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97
4	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	96
5	Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020094.D
Acq On : 23 Apr 2024 04:20
Operator : MA/JU
Sample : P2184-02
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CORE5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.093	19.0	ng/ul	192116000	1	7.705	201774000	20.0
Benzene, 1,4-di...	7.822	20.0	ng/ul	201774000	1	7.705	201774000	20.0
Benzene, 1,2-di...	8.181	26.0	ng/ul	262496000	1	7.705	201774000	20.0
Benzene, 1-brom...	9.152	15.5	ng/ul	1134450	2	10.446	1461900	20.0
m-Chloroaniline	9.475	100.6	ng/ul	7356130	2	10.446	1461900	20.0
Phenol, 2,5-dic...	10.134	11.9	ng/ul	872770	2	10.446	1461900	20.0
Phenol, 2,3-dic...	10.228	47.4	ng/ul	3461070	2	10.446	1461900	20.0
Benzene, 1,3,5-...	10.340	1140.0	ng/ul	83328200	2	10.446	1461900	20.0
Benzene, 1,2,3-...	10.828	305.1	ng/ul	22300900	2	10.446	1461900	20.0
Benzenamine, 2,...	12.728	10.3	ng/ul	887966	3	14.322	1725030	20.0
Phenol, 3,4-dic...	13.340	10.5	ng/ul	908946	3	14.322	1725030	20.0
Benzenamine, 3,...	13.663	3.9	ng/ul	339862	3	14.322	1725030	20.0