

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
 Data File : BM019450.D
 Acq On : 26 Mar 2019 02:27
 Operator : JU/SJ
 Sample : K2027-19
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09E3

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.957	326	335	360	rBV	34257221	54690772	36.29%	10.820%
2	6.787	640	646	657	rBV	150212	243108	0.16%	0.048%
3	6.957	669	675	691	rBV	528099	868985	0.58%	0.172%
4	7.140	699	706	717	rBV	601608	962218	0.64%	0.190%
5	7.498	757	767	780	rBV	23588393	40675746	26.99%	8.047%
6	7.669	780	796	801	rBV2	43933879	102032997	67.70%	20.187%
7	7.998	836	852	857	rBV2	47291170	150702432	100.00%	29.815%
8	8.316	900	906	923	rBV	411705	724261	0.48%	0.143%
9	8.775	977	984	987	rBV	632374	1123086	0.75%	0.222%
10	8.816	987	991	1004	rVB	1730062	2754359	1.83%	0.545%
11	9.092	1031	1038	1045	rBV	158873	255305	0.17%	0.051%
12	9.404	1086	1091	1099	rVB	94859	167179	0.11%	0.033%
13	9.486	1099	1105	1121	rVB	510472	955610	0.63%	0.189%
14	10.016	1188	1195	1209	rBV	774838	1342834	0.89%	0.266%
15	10.269	1227	1238	1244	rBV	33445655	66702278	44.26%	13.197%
16	10.392	1252	1259	1268	rVB	917694	1446174	0.96%	0.286%
17	10.781	1313	1325	1331	rBV	21216466	37224514	24.70%	7.365%
18	11.004	1355	1363	1380	rBV	3884217	6293012	4.18%	1.245%
19	11.216	1391	1399	1414	rBV	816308	1612664	1.07%	0.319%
20	11.710	1470	1483	1497	rVB4	68308	179975	0.12%	0.036%
21	12.427	1591	1605	1616	rBV	764599	1607602	1.07%	0.318%
22	12.845	1671	1676	1692	rVB	179375	292071	0.19%	0.058%
23	13.045	1702	1710	1722	rBV	2727543	4142671	2.75%	0.820%
24	13.680	1810	1818	1823	rBV	1423370	2154508	1.43%	0.426%
25	13.733	1823	1827	1845	rVB	237469	455051	0.30%	0.090%
26	13.951	1857	1864	1879	rBV	1779179	2505853	1.66%	0.496%
27	14.257	1910	1916	1923	rBV2	1197988	1835517	1.22%	0.363%
28	15.257	2079	2086	2094	rBV	2044387	2985913	1.98%	0.591%
29	15.398	2105	2110	2124	rVB	596656	890047	0.59%	0.176%
30	17.015	2379	2385	2396	rBV2	1470391	2023817	1.34%	0.400%
31	17.115	2396	2402	2420	rBV2	2225180	3117199	2.07%	0.617%
32	19.421	2788	2794	2805	rBV	2655655	3538840	2.35%	0.700%
33	21.221	3095	3100	3108	rBV	1806188	2321412	1.54%	0.459%
34	23.291	3445	3452	3466	rBV2	2190354	4003704	2.66%	0.792%

LSC Area Percent Report

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Integration Parameters: LSCINT.P

Integrator: RTE
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Min Area: 0.1 % of largest Peak
Max Peaks: 100
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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35	23.433	3470	3476	3487	rVB2	1284998	2439401	1.62%	0.483%
36	23.915	3554	3558	3568	rVB2	104496	180331	0.12%	0.036%

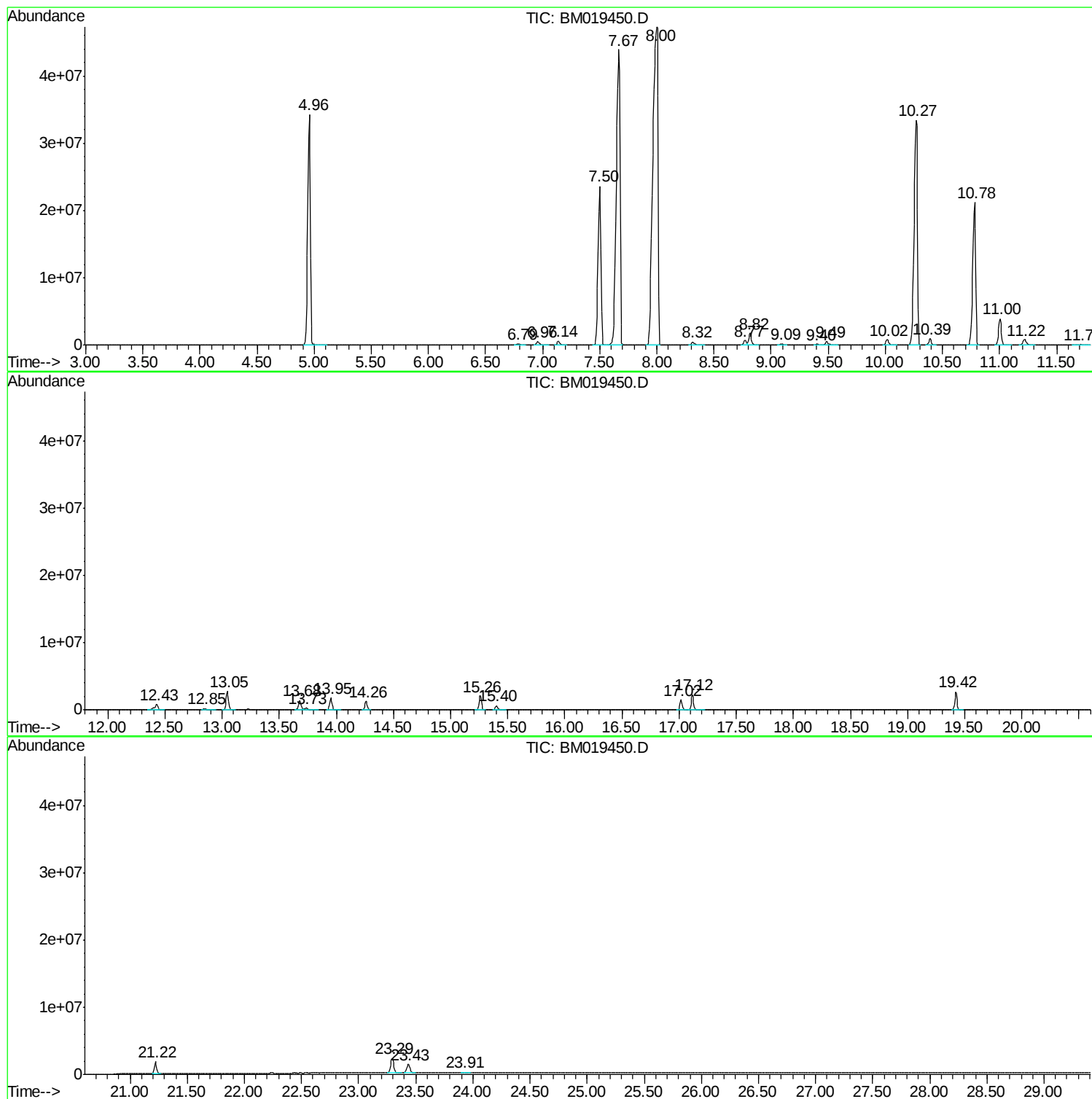
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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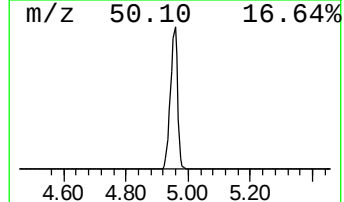
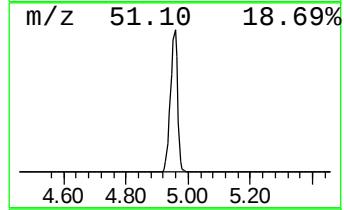
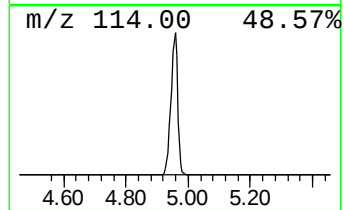
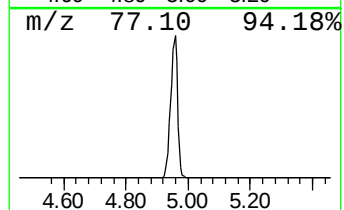
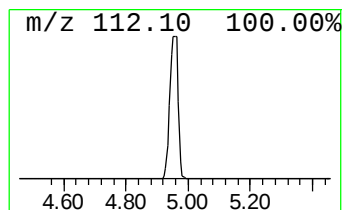
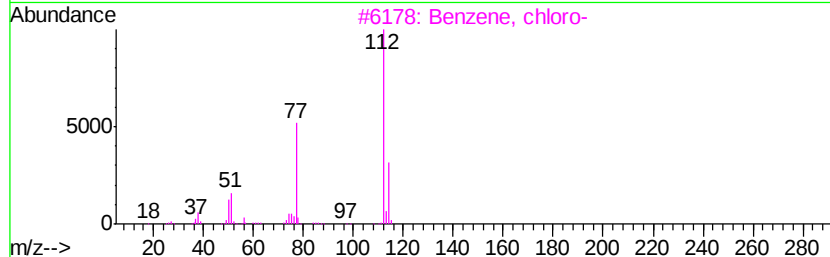
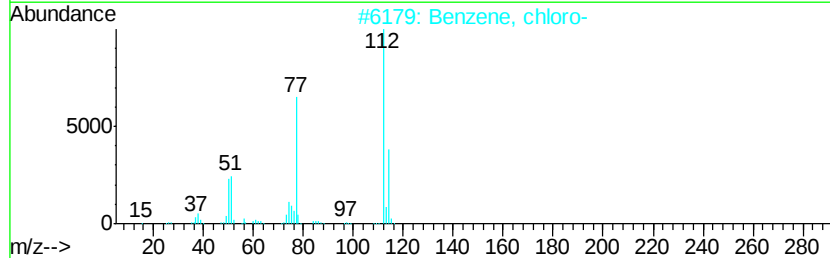
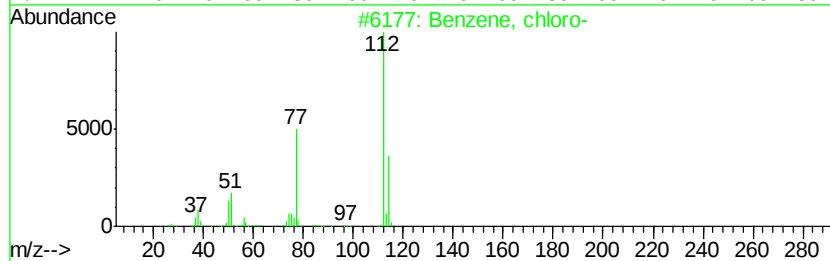
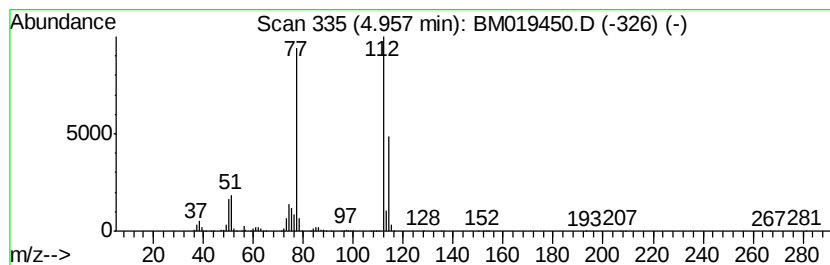
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzene, chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	10.72 ng/ul	54690800	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	94
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		Phenol, 4-fluoro-	112	C6H5FO	000371-41-5	12



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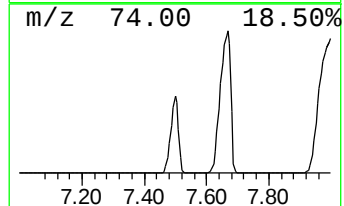
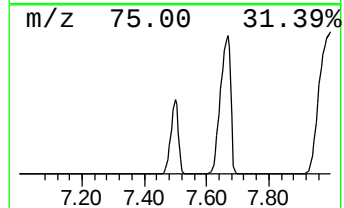
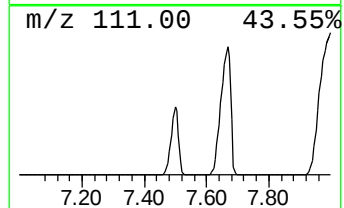
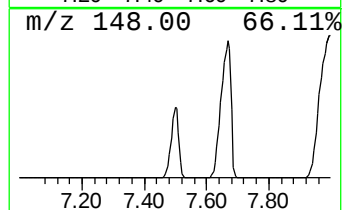
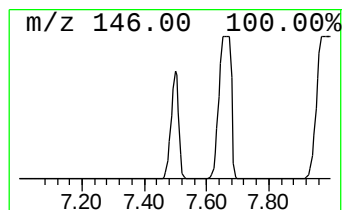
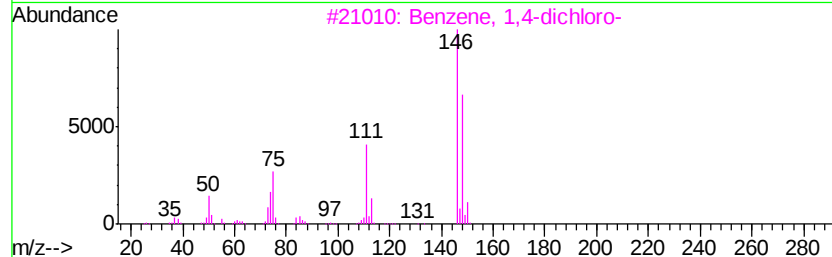
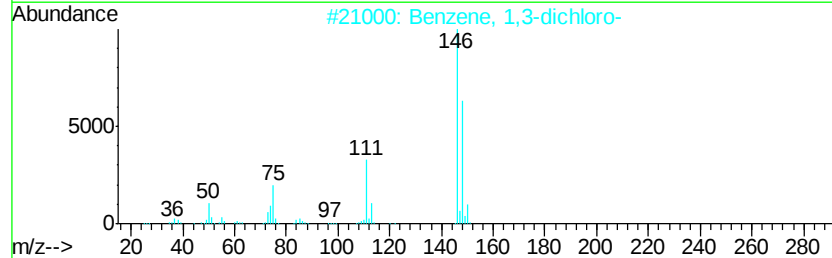
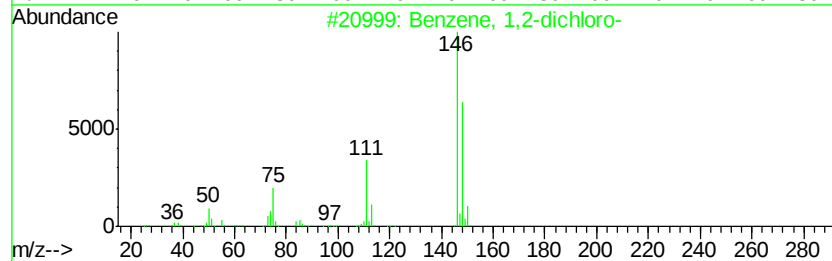
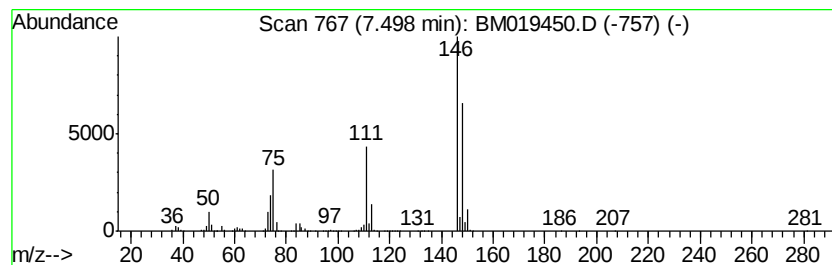
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	7.97 ng/ul	40675700	1,4-Dichlorobenzene-d4	7.61

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



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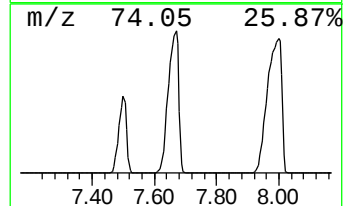
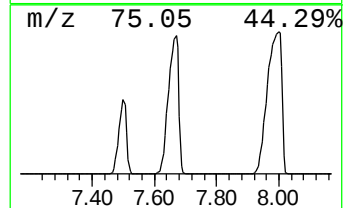
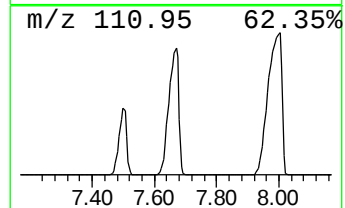
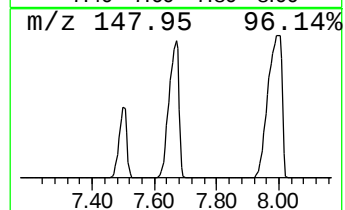
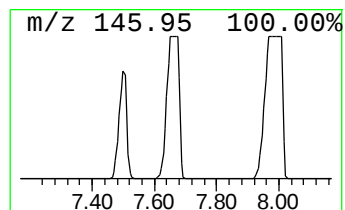
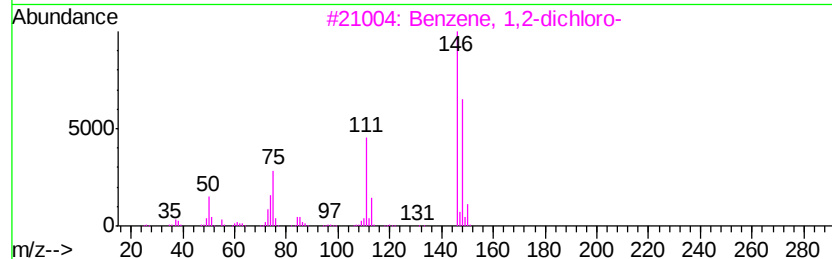
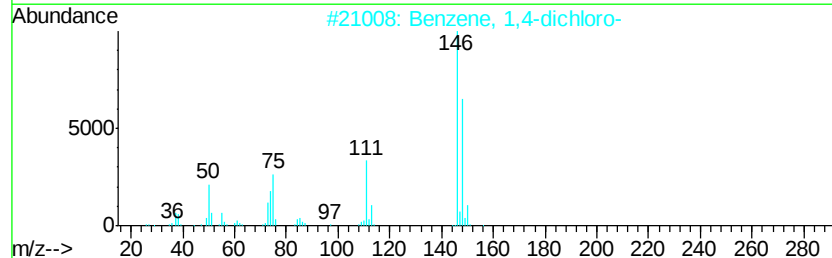
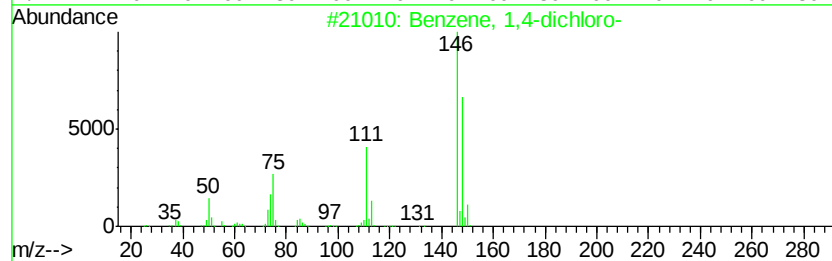
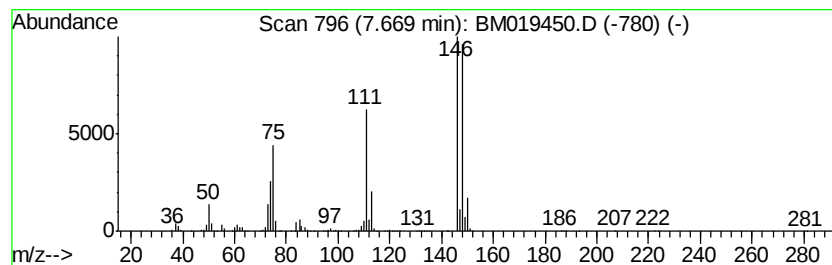
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 3 Benzene, 1,4-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	20.00 ng/ul	102033000	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	91
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	90



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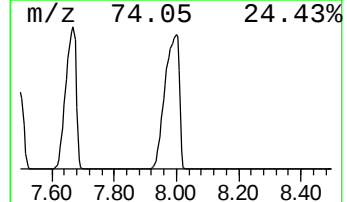
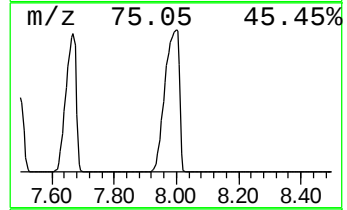
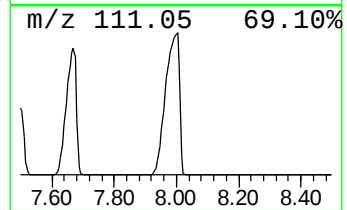
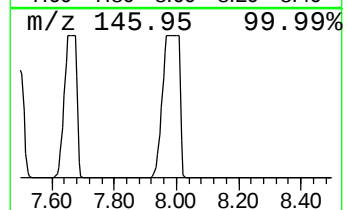
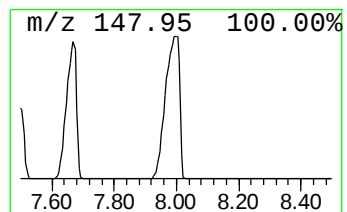
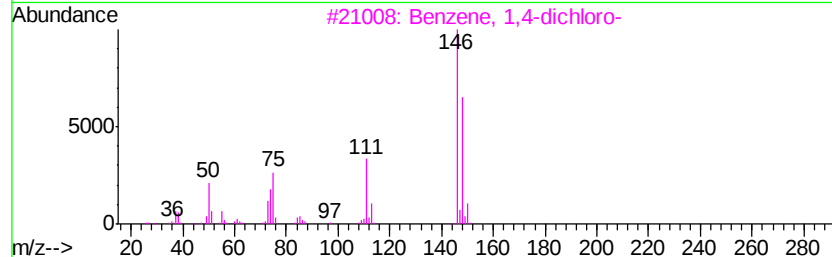
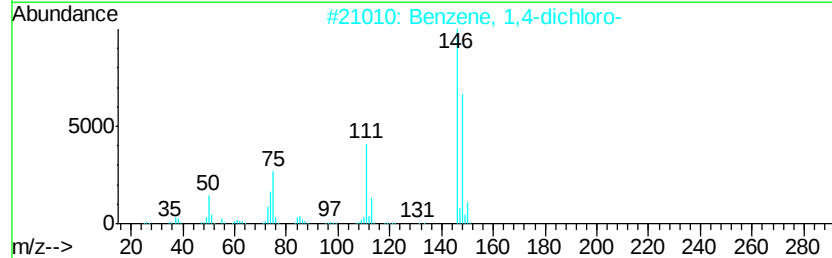
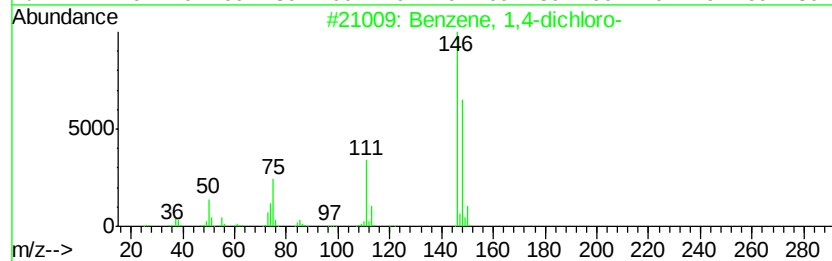
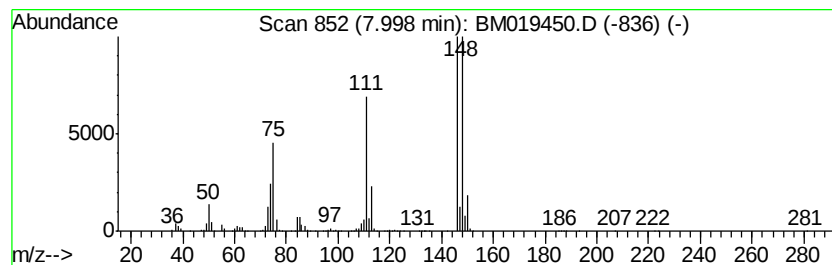
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Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.00	29.54 ng/ul	150702000	1,4-Dichlorobenzene-d4	7.61

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	87



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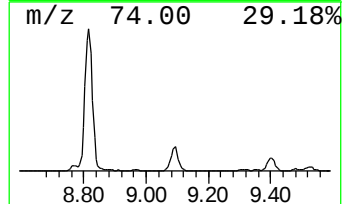
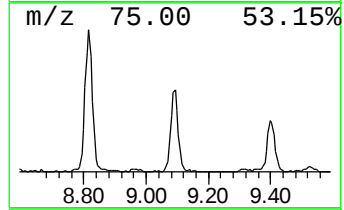
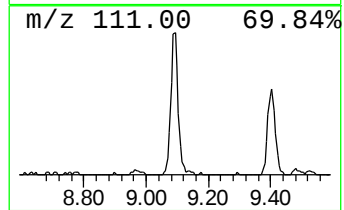
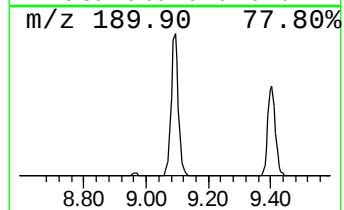
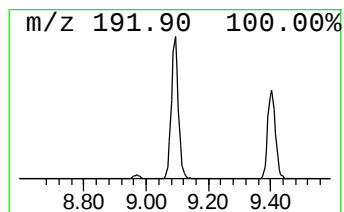
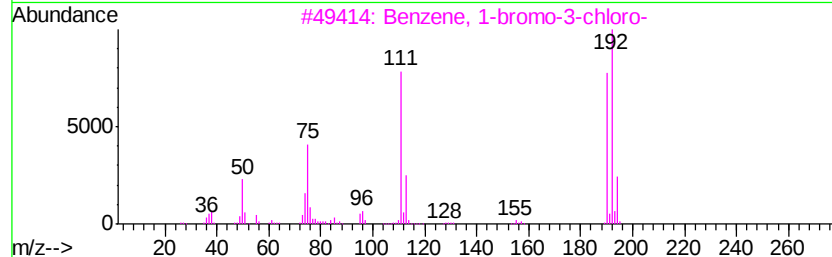
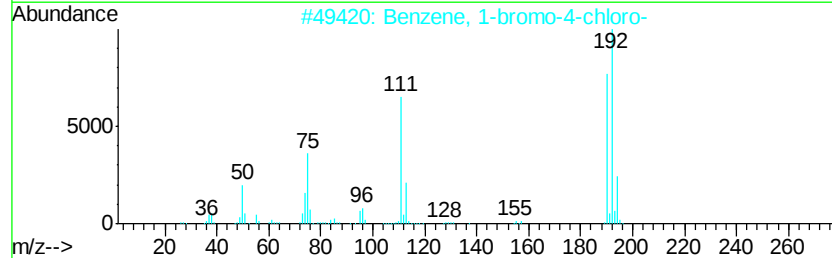
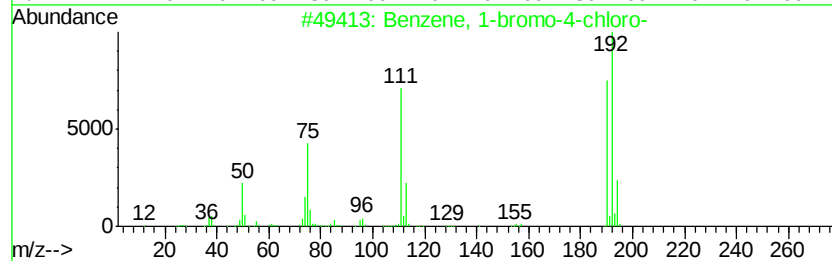
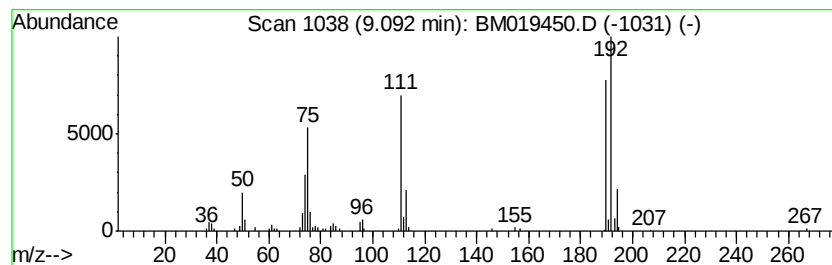
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TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.09	3.53 ng/ul	255305	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
2		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	96
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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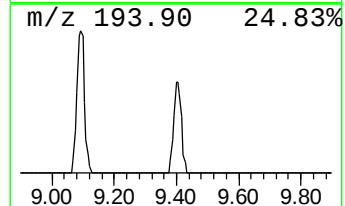
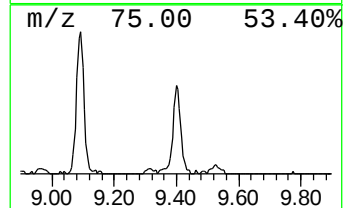
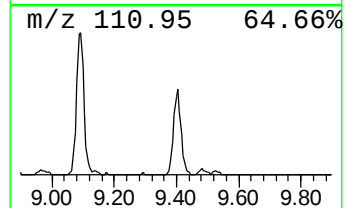
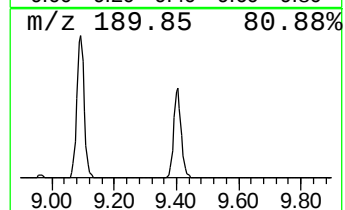
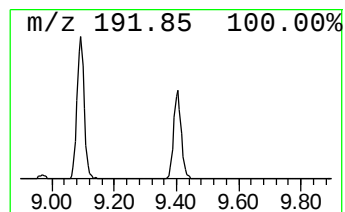
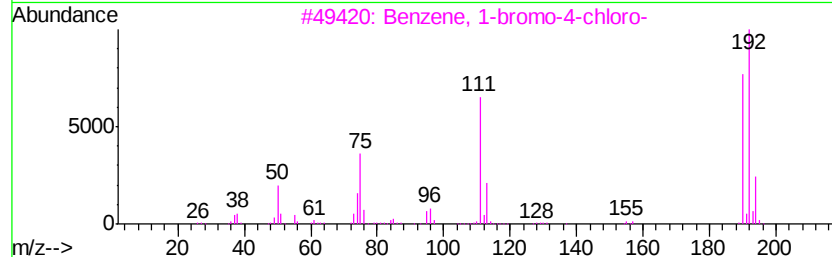
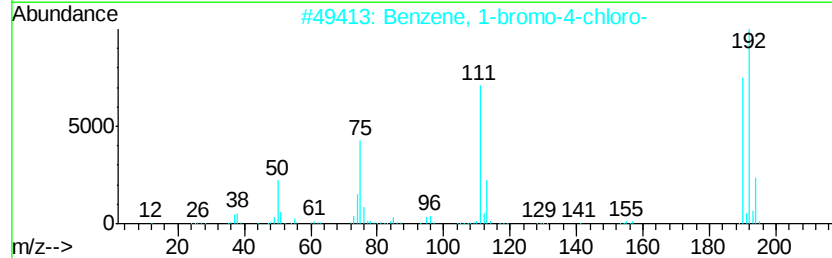
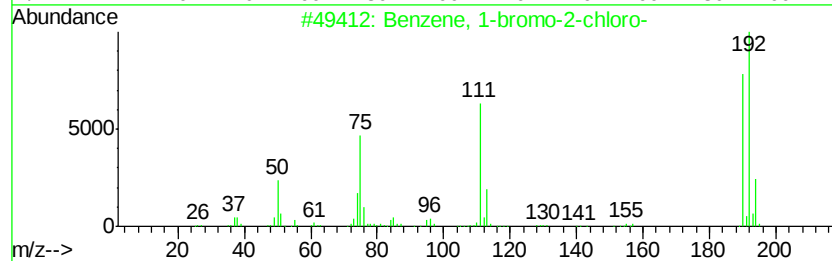
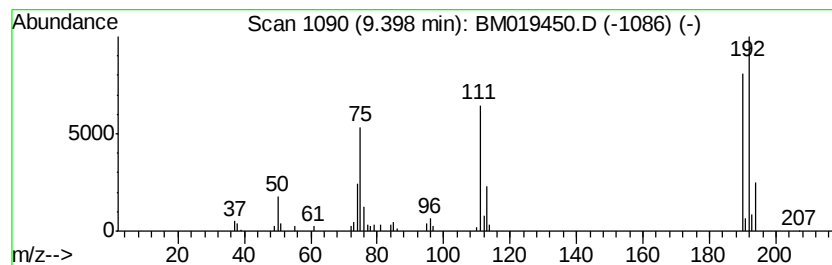
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	2.31 ng/ul	167179	Napthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	97
2		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
4		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	94
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

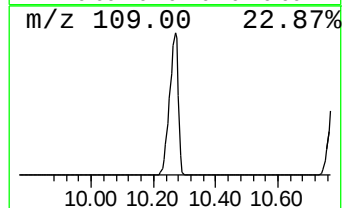
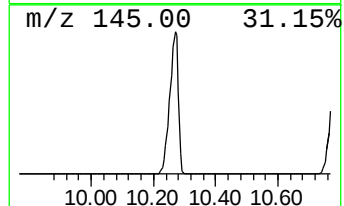
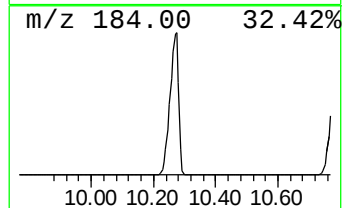
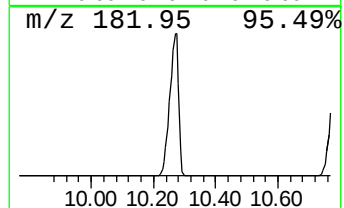
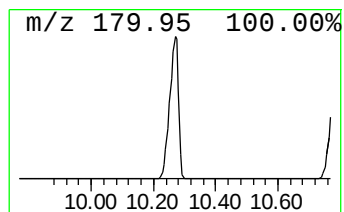
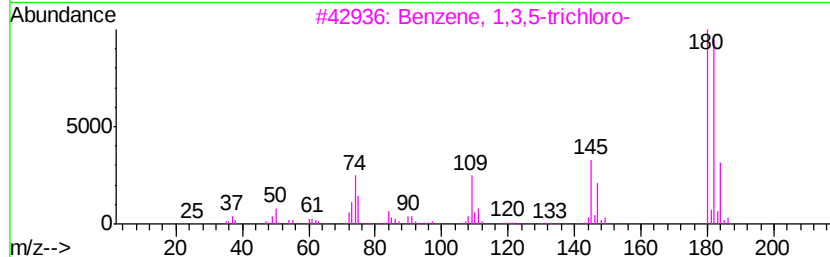
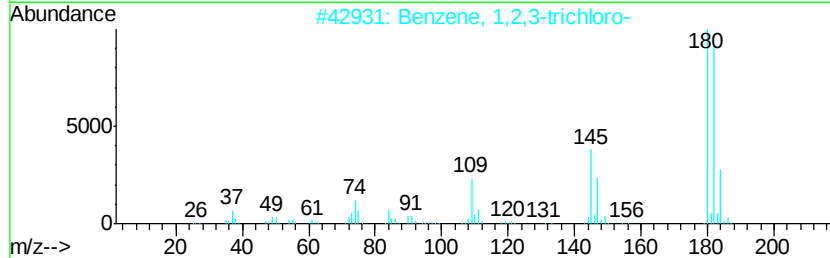
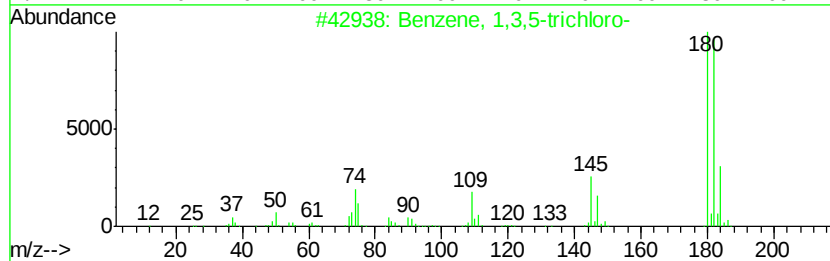
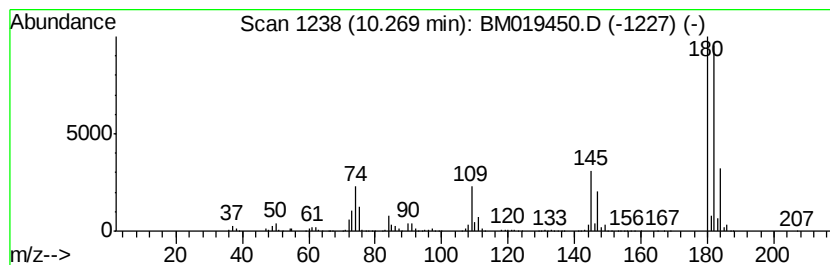
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 1,3,5-trichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.27	922.47 ng/ul	66702300	Napthalene-d8	10.39
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 98
3		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 97
4		Benzene, 1,2,4-trichloro-	180 C6H3Cl3	000120-82-1 95
5		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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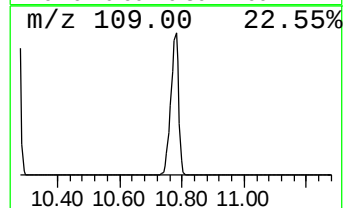
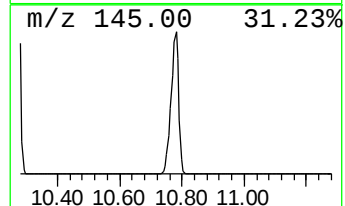
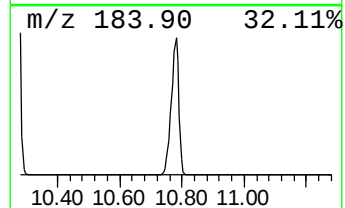
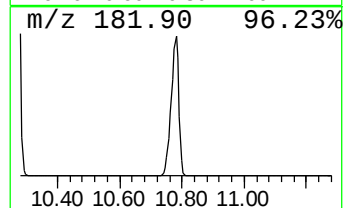
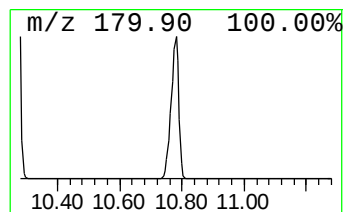
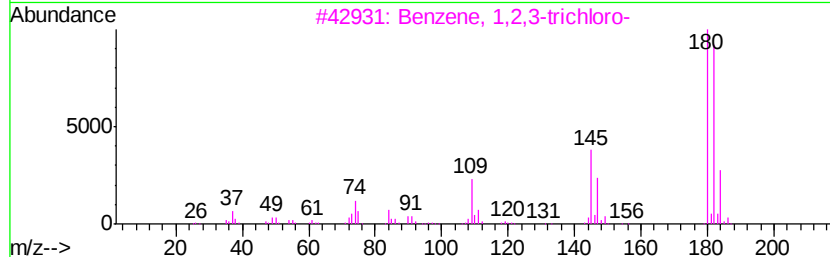
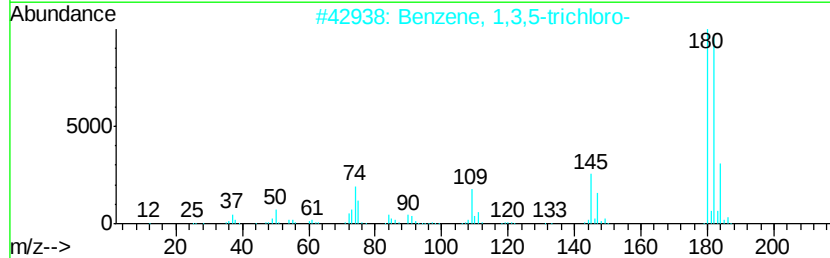
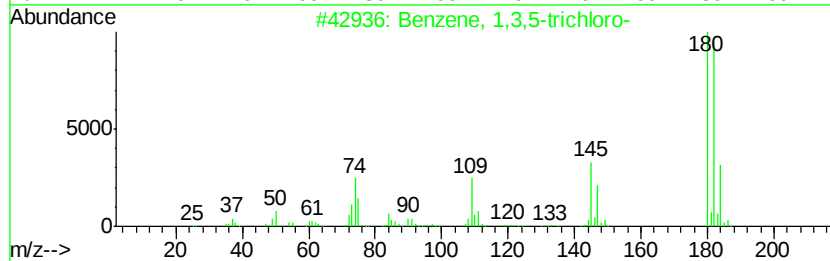
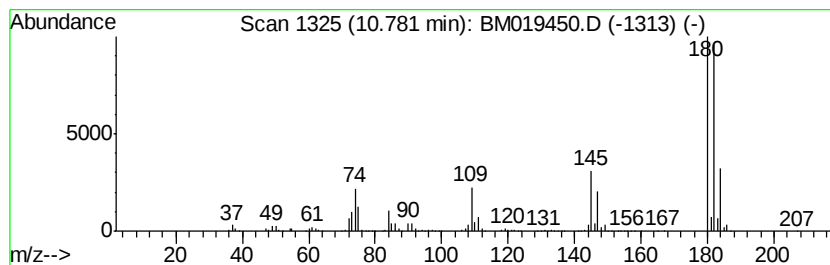
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	514.80 ng/ul	37224500	Napthalene-d8	10.39

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	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 M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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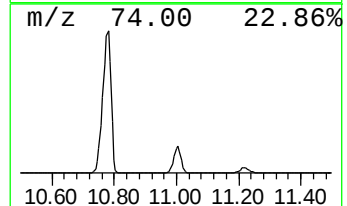
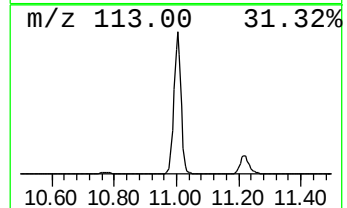
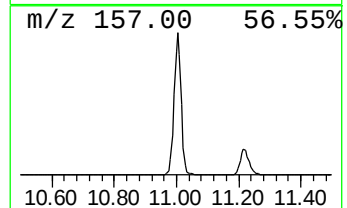
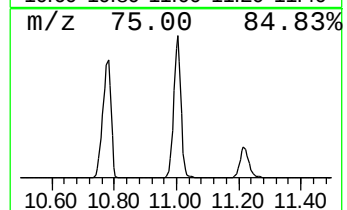
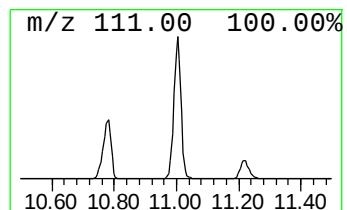
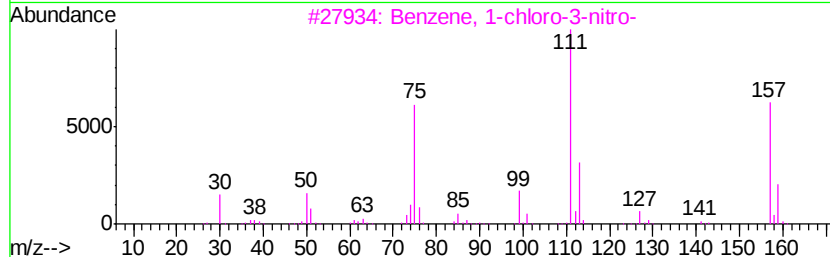
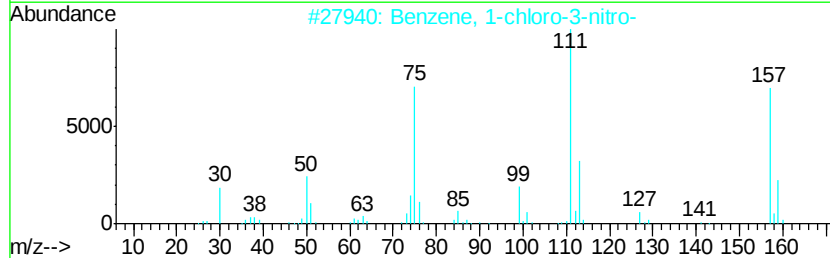
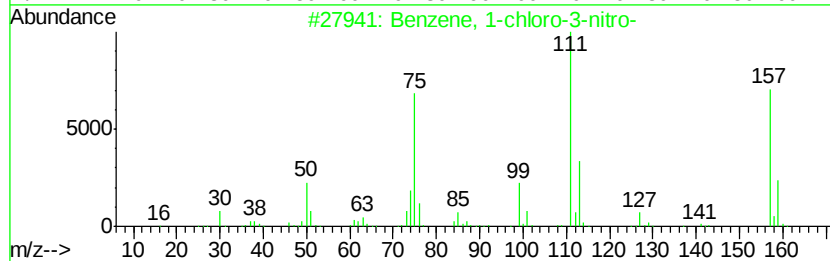
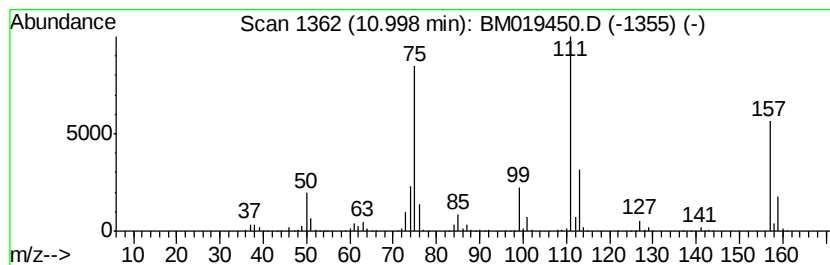
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.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.
11.00	87.03 ng/ul	6293010	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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-3-nitro-	157	C6H4ClNO2	000121-73-3	95
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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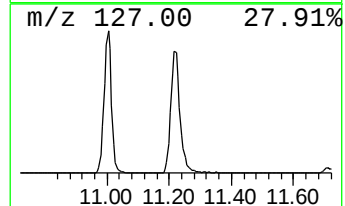
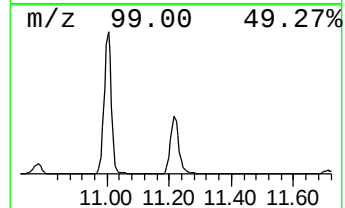
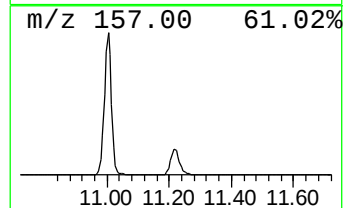
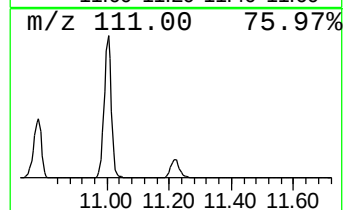
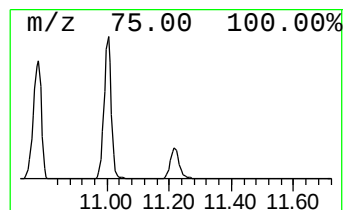
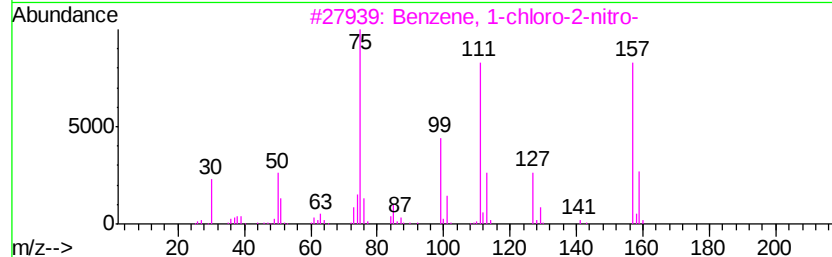
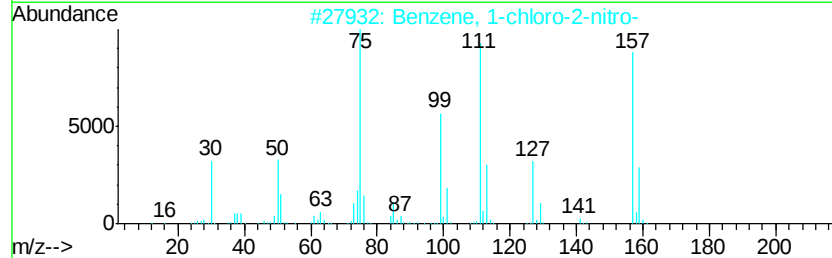
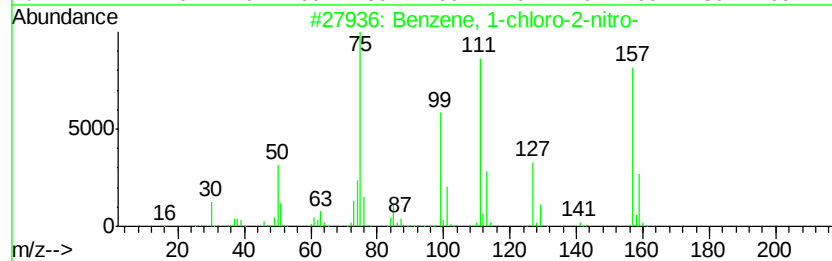
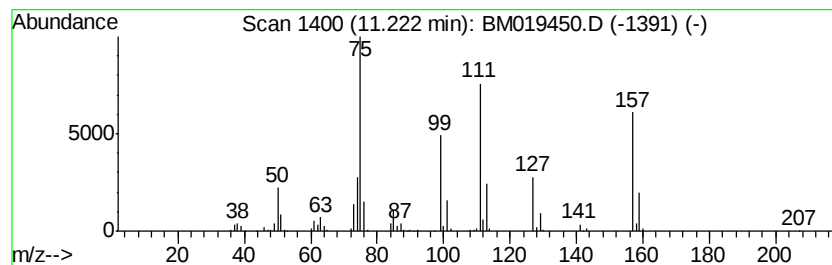
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-chloro-4-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	22.30 ng/ul	1612660	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	99
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

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ClientSampleId :
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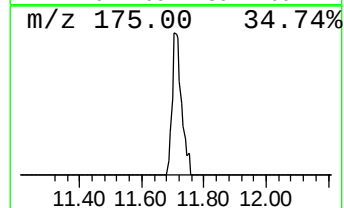
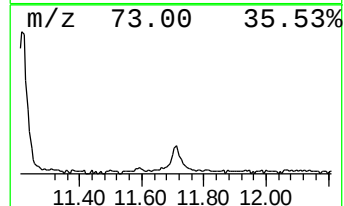
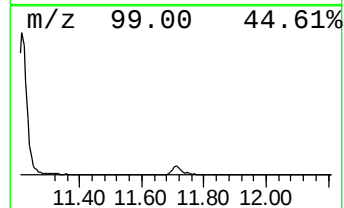
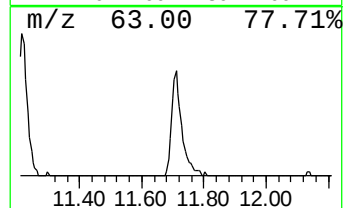
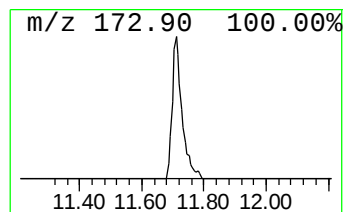
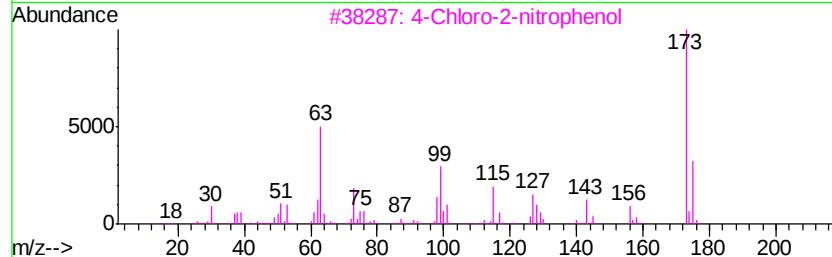
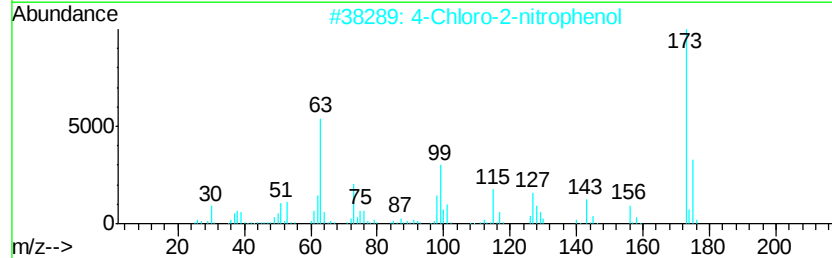
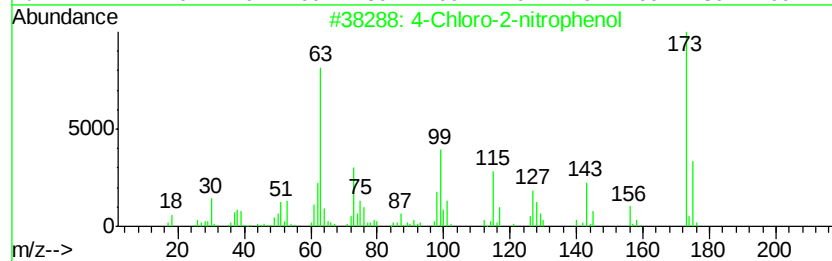
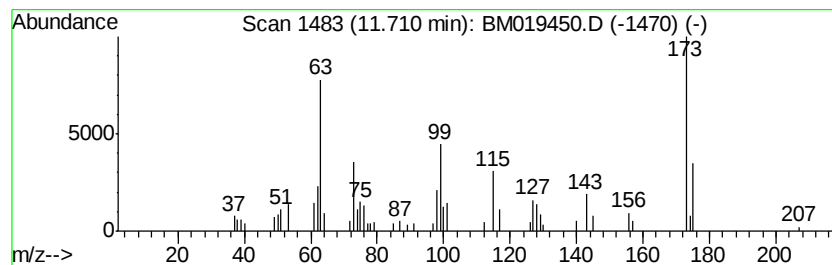
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 4-Chloro-2-nitrophenol Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.71	2.49 ng/ul	179975	Naphthalene-d8	10.39

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	96
3			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	94
4			Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	93
5			Phenol, 2-chloro-4-nitro-	173	C6H4ClNO3	000619-08-9	68



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
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Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

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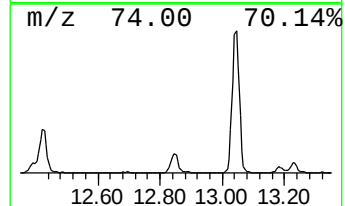
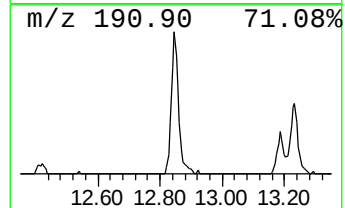
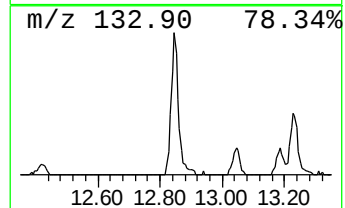
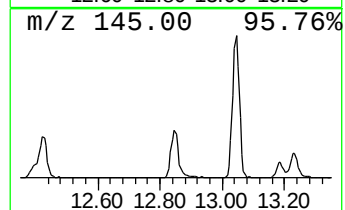
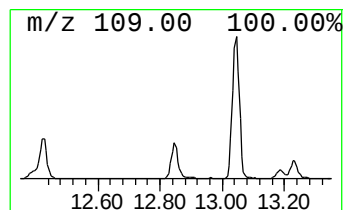
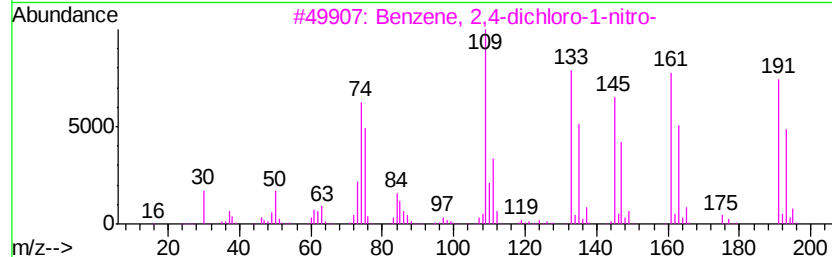
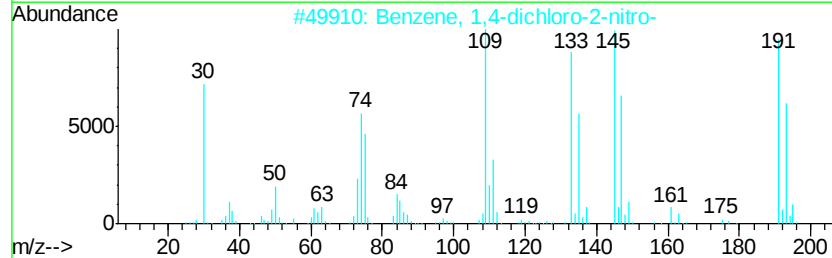
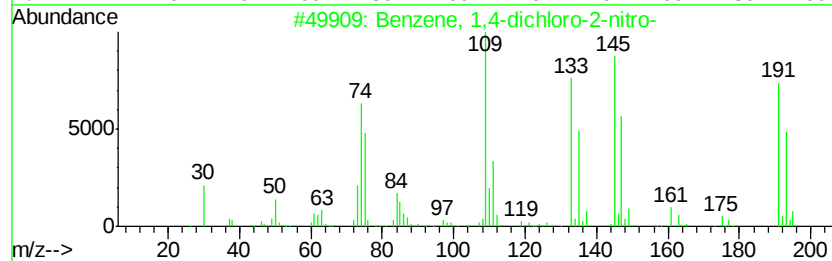
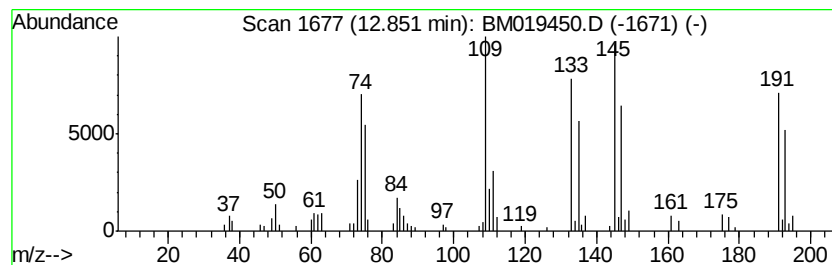
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Benzene, 1,4-dichloro-2-nitro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.85	3.18 ng/ul	292071	Acenaphthene-d10	14.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	99
2			Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	96
3			Benzene, 2,4-dichloro-1-nitro-	191	C6H3Cl2NO2	000611-06-3	94
4			Benzene, 2,4-dichloro-1-nitro-	191	C6H3Cl2NO2	000611-06-3	94
5			Benzene, 1,2-dichloro-3-nitro-	191	C6H3Cl2NO2	003209-22-1	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032519\
Data File : BM019450.D
Acq On : 26 Mar 2019 02:27
Operator : JU/SJ
Sample : K2027-19
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C09E3

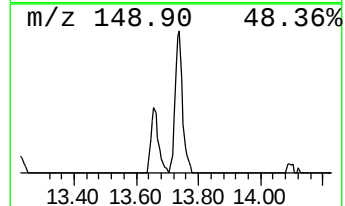
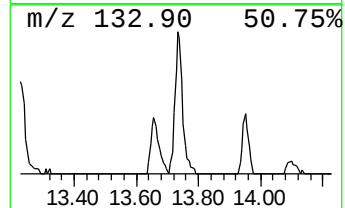
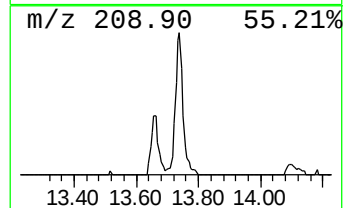
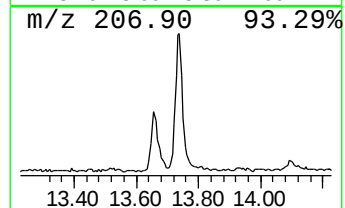
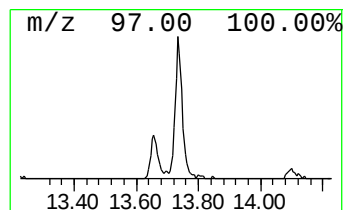
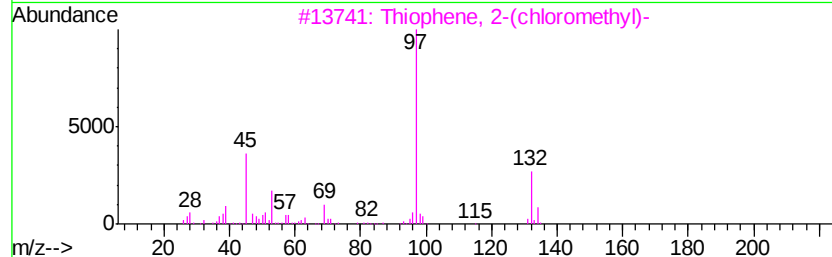
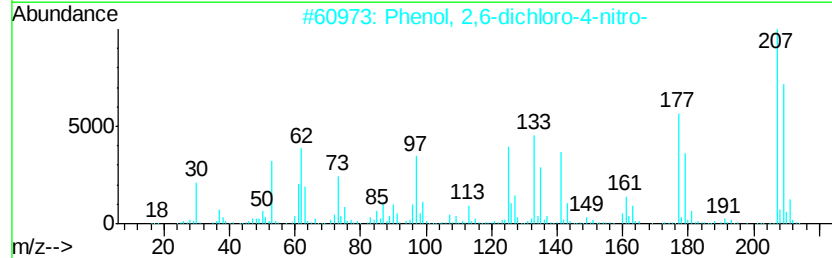
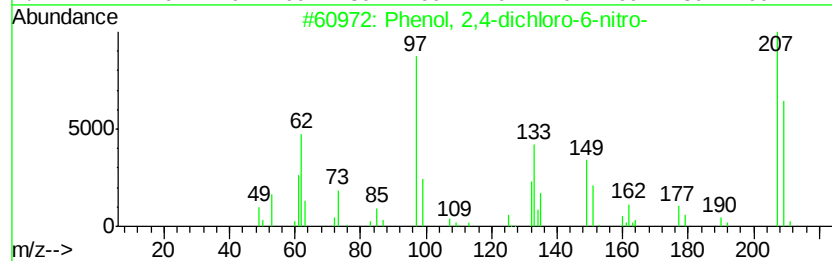
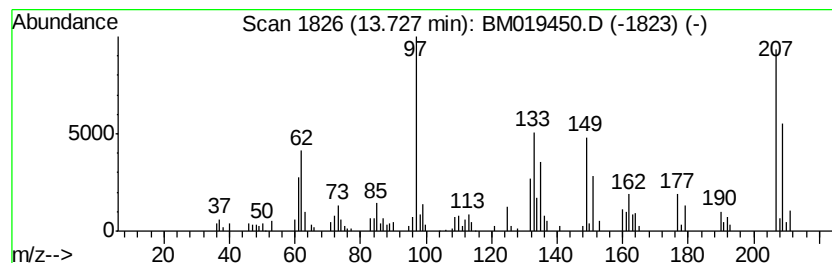
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Phenol, 2,4-dichloro-6-nitro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.73	4.96 ng/ul	455051	Acenaphthene-d10	14.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,4-dichloro-6-nitro-	207	C6H3Cl2NO3	000609-89-2	94
2		Phenol, 2,6-dichloro-4-nitro-	207	C6H3Cl2NO3	000618-80-4	60
3		Thiophene, 2-(chloromethyl)-	132	C5H5ClS	000765-50-4	11
4		4-Hydroxyphenyl pyrrolidinyl thione	207	C11H13NOS	084783-02-8	11
5		2-Nitro-4-(trifluoromethyl)phenol	207	C7H4F3NO3	000400-99-7	10



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032519\
 Data File : BM019450.D
 Acq On : 26 Mar 2019 02:27
 Operator : JU/SJ
 Sample : K2027-19
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C09E3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	4.96	10.7	ng/ul	54690800	1	7.61	102033000	20.0
Benzene, 1,2-dich...	7.50	8.0	ng/ul	40675700	1	7.61	102033000	20.0
Benzene, 1,4-dich...	7.67	20.0	ng/ul	102033000	1	7.61	102033000	20.0
Benzene, 1,3-dich...	8.00	29.5	ng/ul	150702000	1	7.61	102033000	20.0
Benzene, 1-bromo-...	9.09	3.5	ng/ul	255305	2	10.39	1446170	20.0
Benzene, 1-bromo-...	9.40	2.3	ng/ul	167179	2	10.39	1446170	20.0
Benzene, 1,3,5-tr...	10.27	922.5	ng/ul	66702300	2	10.39	1446170	20.0
Benzene, 1,2,3-tr...	10.78	514.8	ng/ul	37224500	2	10.39	1446170	20.0
Benzene, 1-chloro...	11.00	87.0	ng/ul	6293010	2	10.39	1446170	20.0
Benzene, 1-chloro...	11.22	22.3	ng/ul	1612660	2	10.39	1446170	20.0
4-Chloro-2-nitrop...	11.71	2.5	ng/ul	179975	2	10.39	1446170	20.0
Benzene, 1,4-dich...	12.85	3.2	ng/ul	292071	3	14.26	1835520	20.0
Phenol, 2,4-dichl...	13.73	5.0	ng/ul	455051	3	14.26	1835520	20.0