

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025816.D
 Acq On : 27 Mar 2020 00:55
 Operator : CG/JU
 Sample : L1968-12
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CODE3

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-BM031420MA.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	4.963	341	353	358	rBV3	64050724	159933387	43.09%	12.509%
2	6.269	565	575	584	rVB	325130	530787	0.14%	0.042%
3	6.928	679	687	697	rBV	531954	1025377	0.28%	0.080%
4	7.110	710	718	728	rBV	668037	1184376	0.32%	0.093%
5	7.492	769	783	793	rBV2	50337248	129412065	34.87%	10.122%
6	7.675	793	814	818	rBV5	72751790	273136720	73.59%	21.362%
7	8.016	849	872	876	rBV6	74034608	371148511	100.00%	29.028%
8	8.275	910	916	925	rVB	552608	863281	0.23%	0.068%
9	8.716	984	991	994	rBV	855349	1484402	0.40%	0.116%
10	8.763	994	999	1009	rVB	3813168	6026008	1.62%	0.471%
11	9.045	1041	1047	1057	rVB	611090	946564	0.26%	0.074%
12	9.357	1095	1100	1106	rVV	467262	785816	0.21%	0.061%
13	9.428	1106	1112	1121	rVV	731943	1316016	0.35%	0.103%
14	9.963	1195	1203	1213	rBV	1028545	1749145	0.47%	0.137%
15	10.251	1236	1252	1256	rBV2	59834099	152250272	41.02%	11.908%
16	10.351	1263	1269	1281	rVB	1080704	1637487	0.44%	0.128%
17	10.751	1324	1337	1342	rBV	43795157	94403693	25.44%	7.383%
18	10.945	1360	1370	1381	rVB	9388423	15976153	4.30%	1.250%
19	11.145	1396	1404	1417	rVB	2566405	4225953	1.14%	0.331%
20	11.627	1476	1486	1502	rVB4	281697	591280	0.16%	0.046%
21	12.322	1598	1604	1607	rBV	552576	980622	0.26%	0.077%
22	12.380	1610	1614	1622	rVB	1850957	2777160	0.75%	0.217%
23	12.769	1674	1680	1686	rBV	580209	856858	0.23%	0.067%
24	12.998	1711	1719	1731	rBV	7470160	10860419	2.93%	0.849%
25	13.157	1741	1746	1752	rVB	282792	409147	0.11%	0.032%
26	13.592	1814	1820	1822	rBV	882276	1310457	0.35%	0.102%
27	13.621	1822	1825	1829	rVV	2057305	2776355	0.75%	0.217%
28	13.669	1829	1833	1841	rVB	728701	1050927	0.28%	0.082%
29	13.892	1859	1871	1878	rBV	2419552	3493420	0.94%	0.273%
30	14.204	1916	1924	1938	rVB	1522678	2319977	0.63%	0.181%
31	14.692	2002	2007	2013	rBV2	287001	388466	0.10%	0.030%
32	15.204	2088	2094	2099	rBV	3172351	4444934	1.20%	0.348%
33	15.327	2109	2115	2125	rVB	1046863	1465177	0.39%	0.115%
34	16.951	2385	2391	2397	rBV	1860502	2586624	0.70%	0.202%

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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Title : SVOA CALIBRATION

35	17.051	2401	2408	2415	rBV	3465870	4791197	1.29%	0.375%
36	19.345	2792	2798	2805	rBV	4295694	5679799	1.53%	0.444%
37	20.886	3056	3060	3067	rBV2	336800	481946	0.13%	0.038%
38	21.139	3098	3103	3110	rBV	2570344	3209076	0.86%	0.251%
39	23.156	3439	3446	3453	rBV	3798536	6583726	1.77%	0.515%
40	23.291	3462	3469	3479	rVB	1995069	3487987	0.94%	0.273%

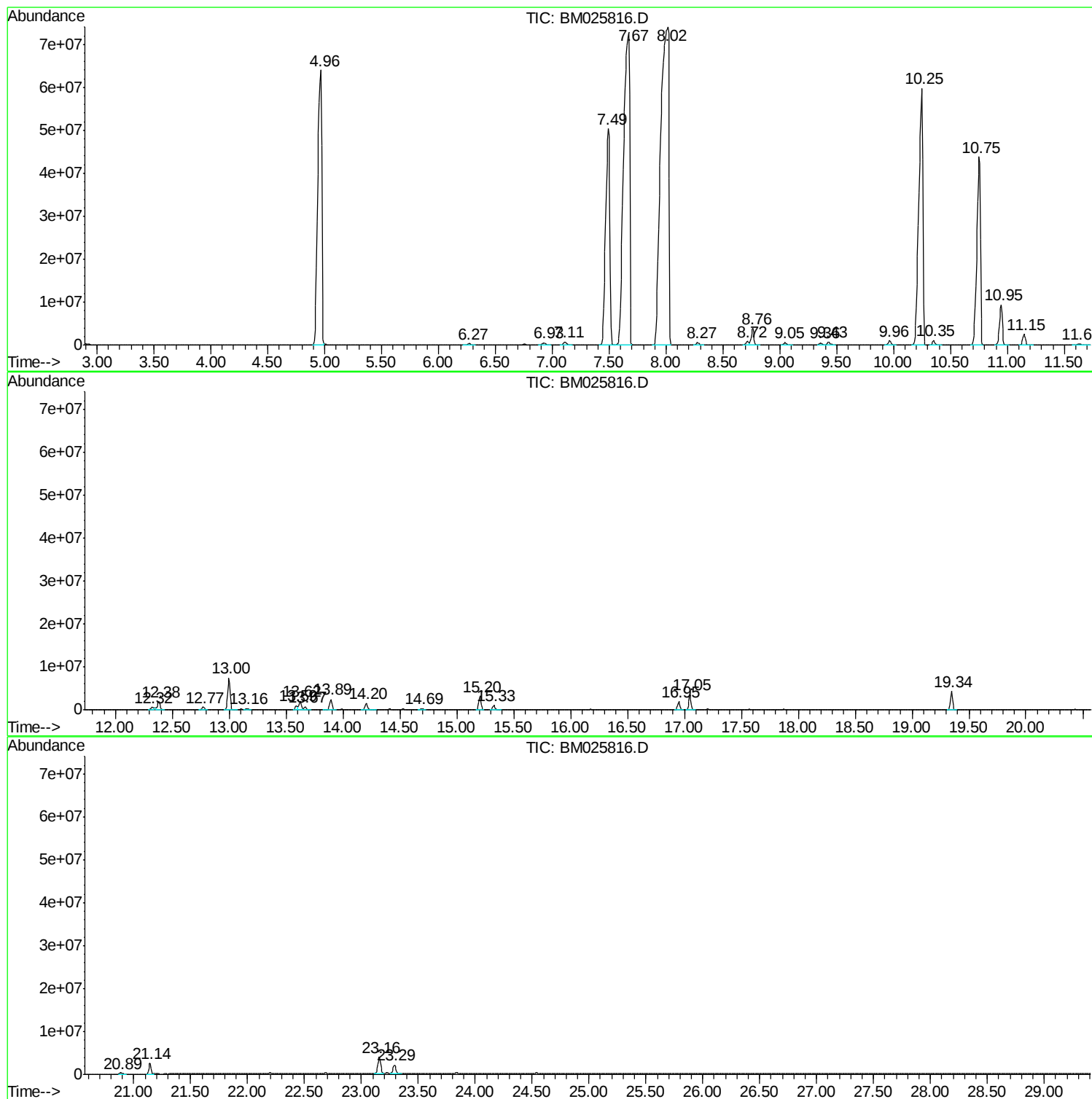
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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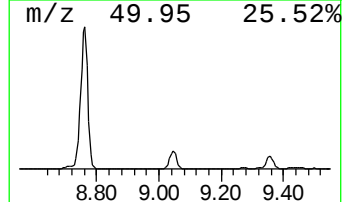
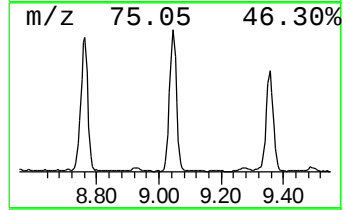
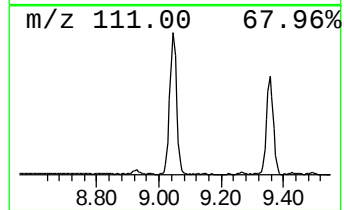
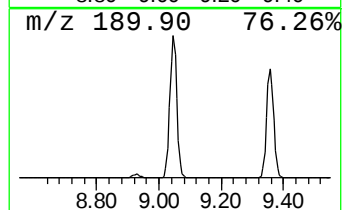
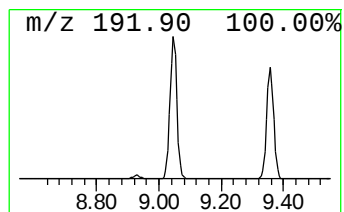
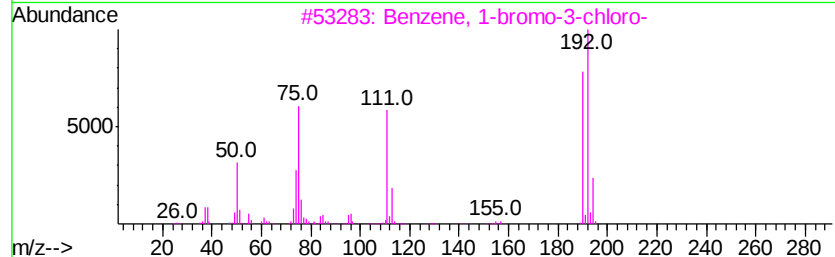
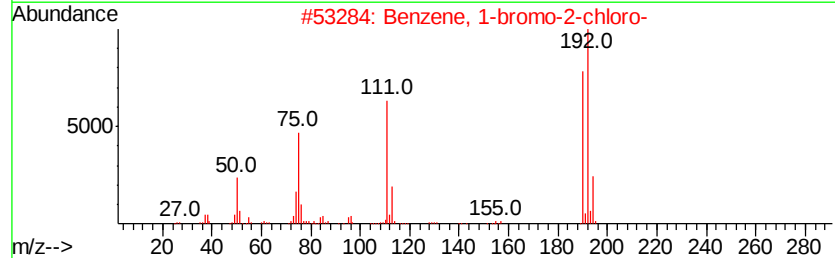
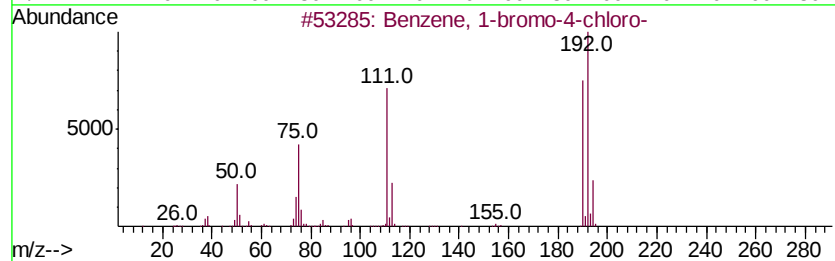
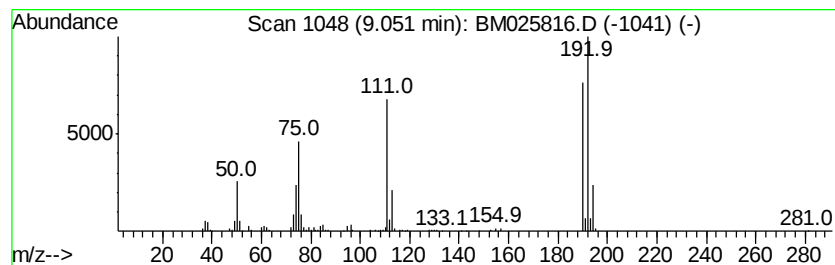
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
9.05	11.56 ng/ul	946564	Napthalene-d8	10.35

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-2-chloro-	190	C6H4BrCl	000694-80-4	98
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
5		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97



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Instrument :
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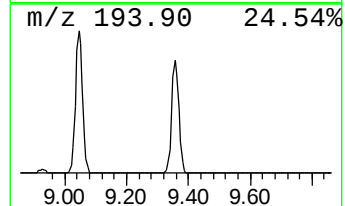
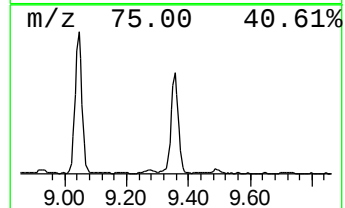
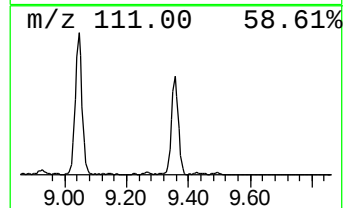
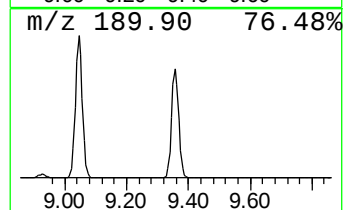
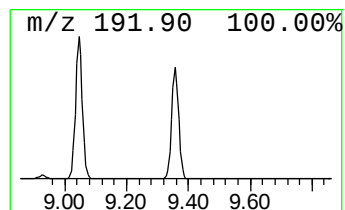
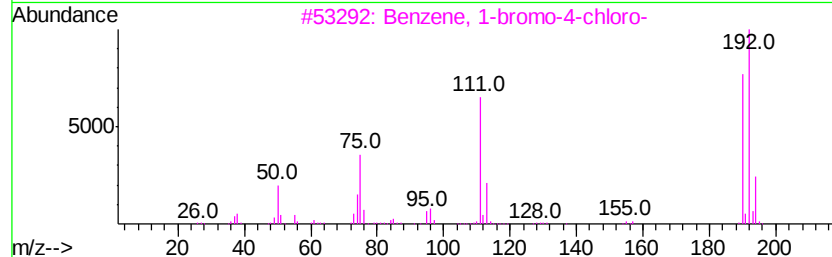
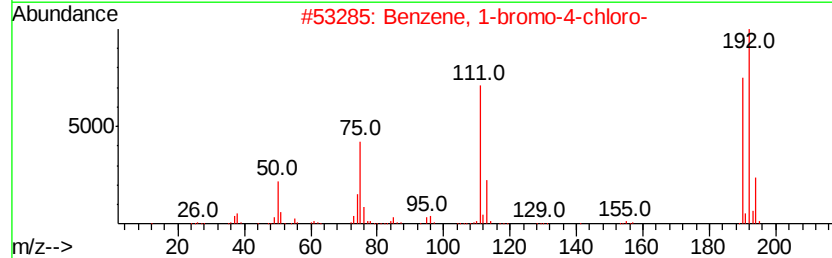
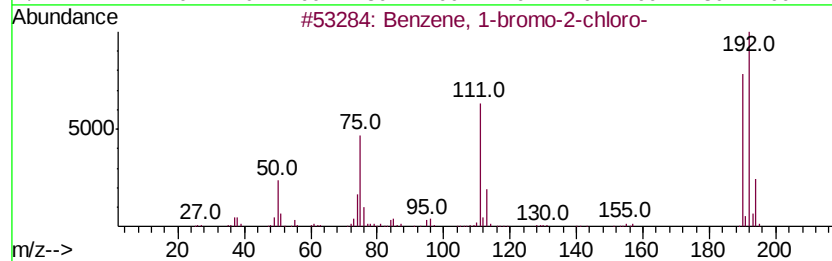
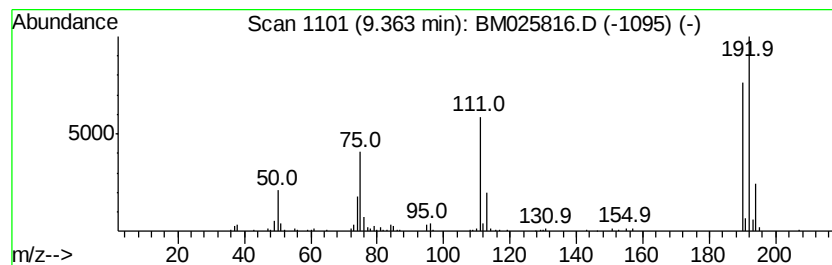
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	9.60 ng/ul	785816	Napthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
4		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	95
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



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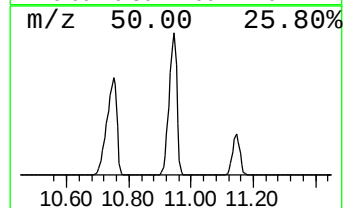
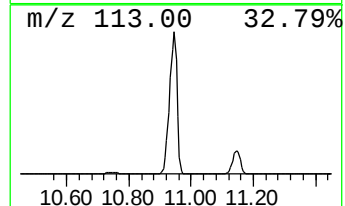
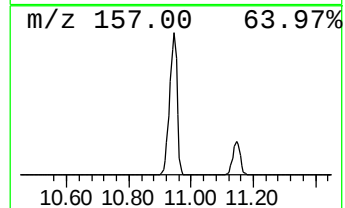
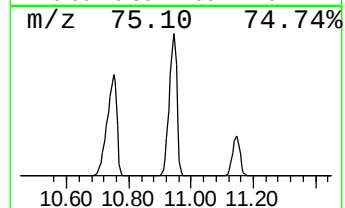
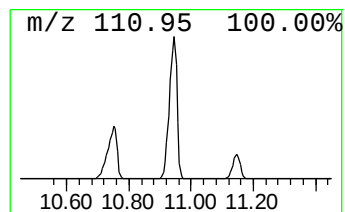
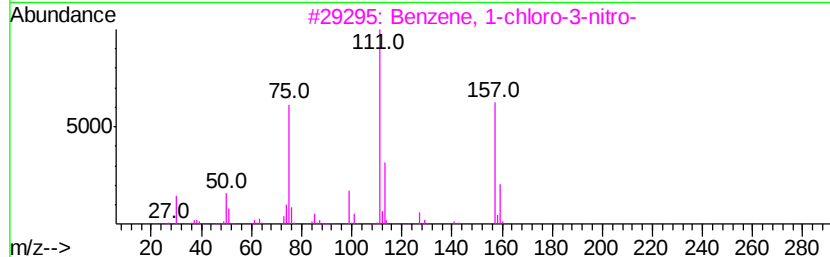
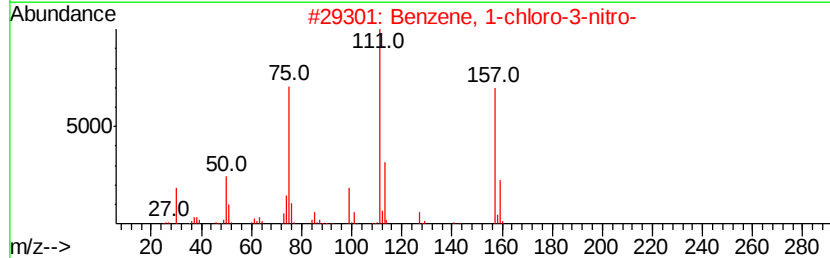
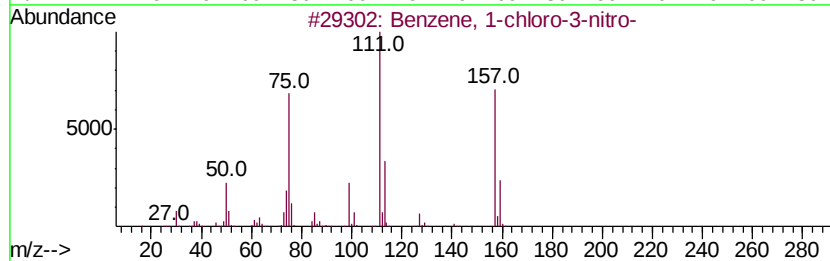
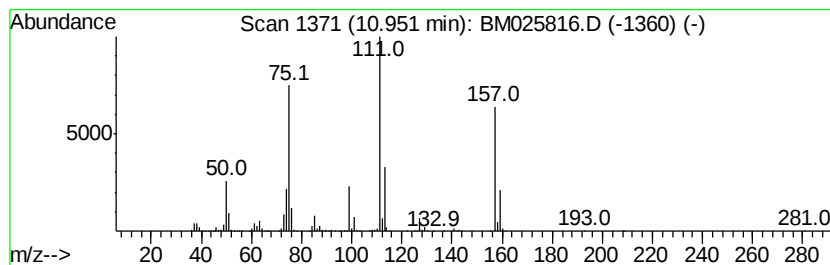
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
10.95	195.13 ng/ul	15976200	Naphthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	98
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91



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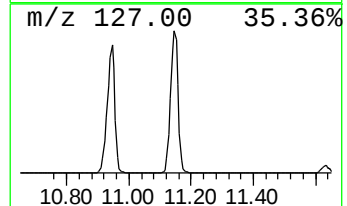
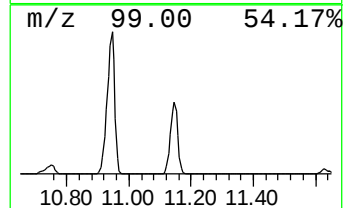
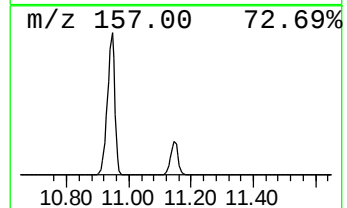
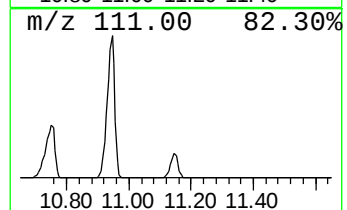
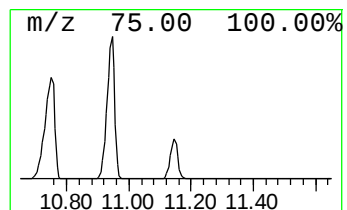
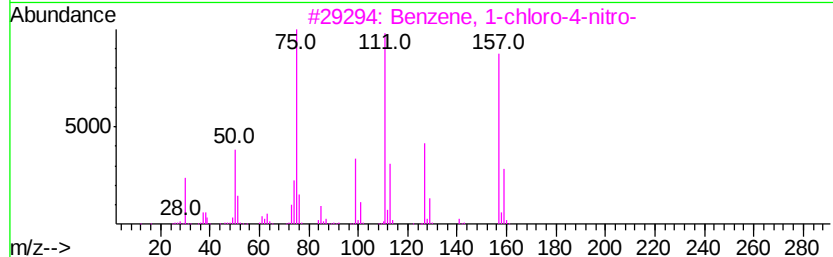
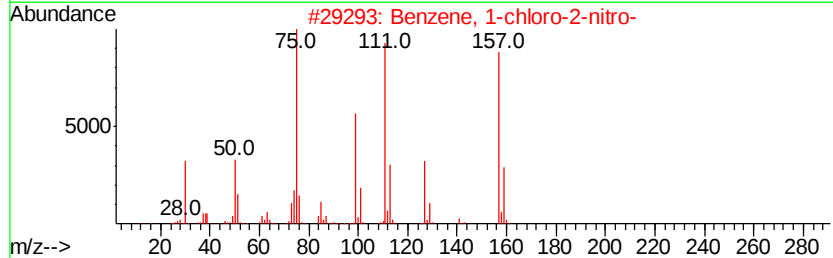
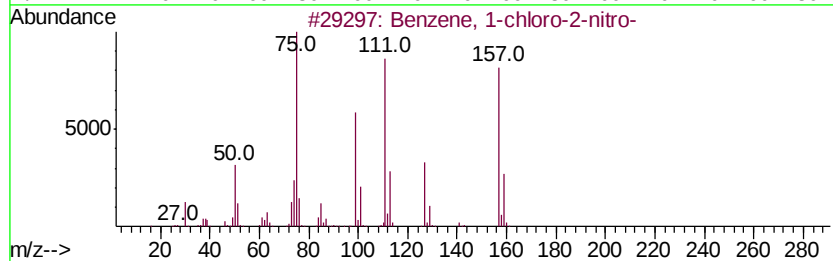
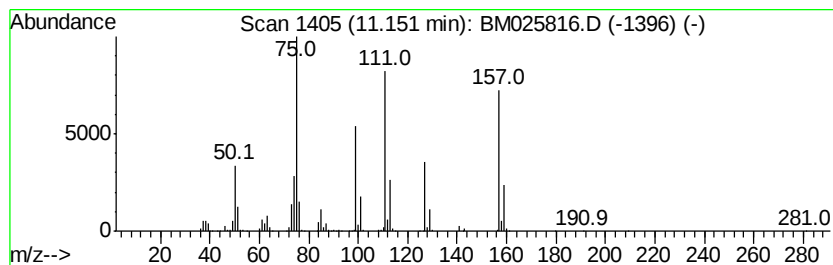
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	51.62 ng/ul	4225950	Napthalene-d8	10.35

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-4-nitro-	157	C6H4ClNO2	000100-00-5	98
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97



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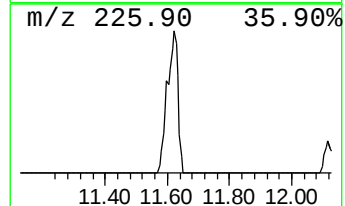
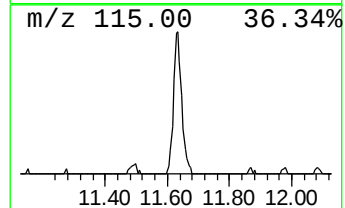
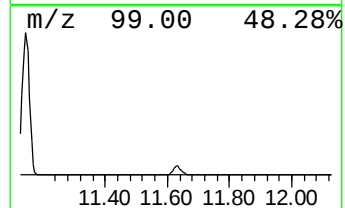
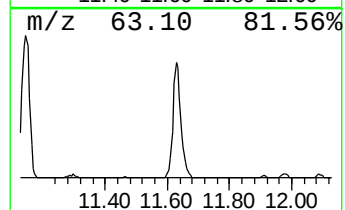
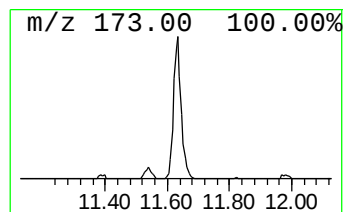
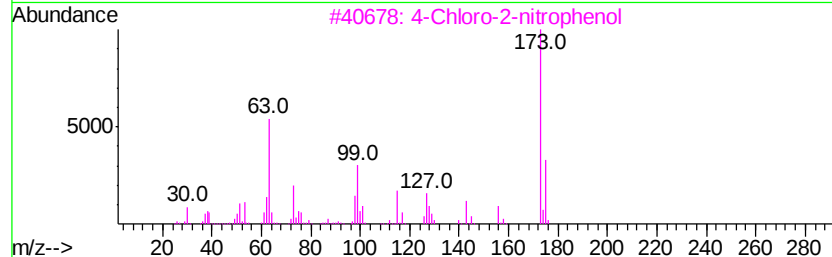
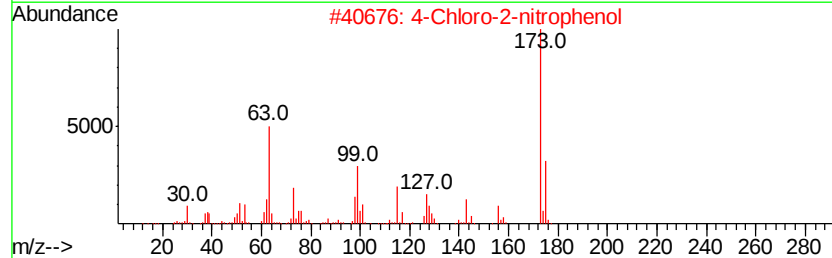
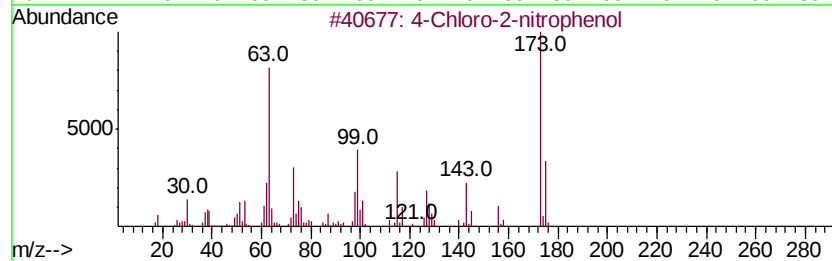
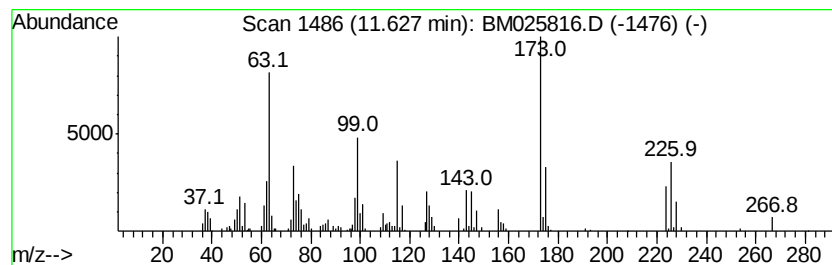
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 4-Chloro-2-nitrophenol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.63	7.22 ng/ul	591280	Naphthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	94
3		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	93
4		Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	64
5		3-Chloro-4-nitrophenol	173	C6H4ClNO3	000491-11-2	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025816.D
Acq On : 27 Mar 2020 00:55
Operator : CG/JU
Sample : L1968-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

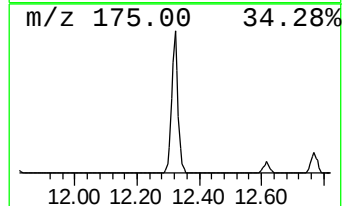
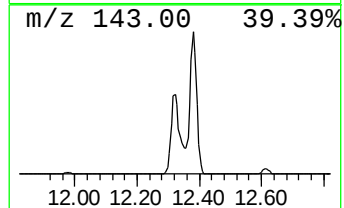
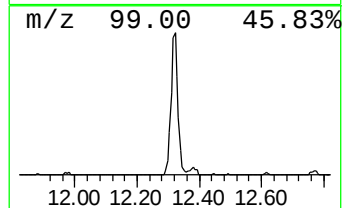
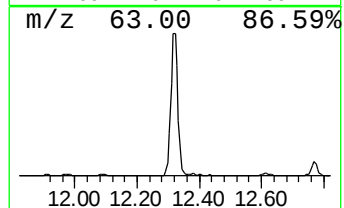
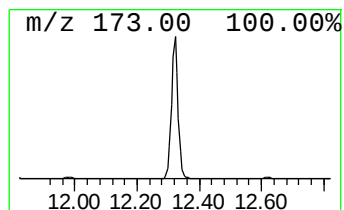
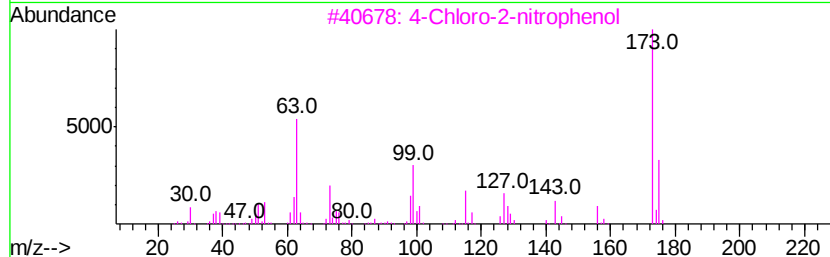
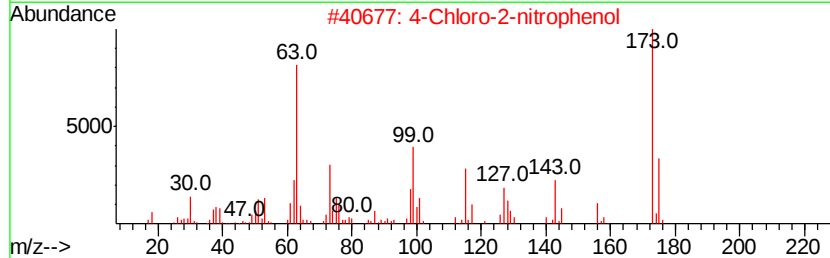
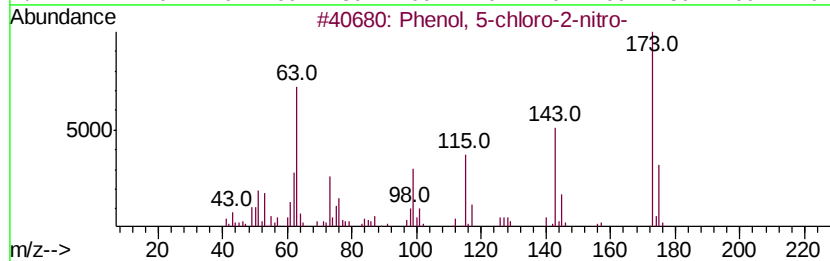
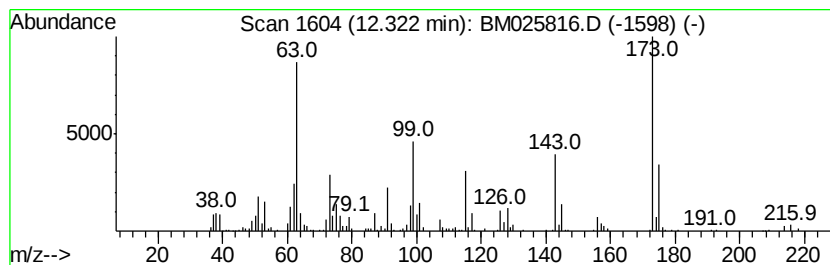
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Phenol, 5-chloro-2-nitro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.32	8.45 ng/ul	980622	Acenaphthene-d10	14.20	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	91
2	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	90
3	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	87
4	Propane, 1,1,2-trichloro-	146	C3H5Cl3	000598-77-6	47
5	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025816.D
Acq On : 27 Mar 2020 00:55
Operator : CG/JU
Sample : L1968-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CODE3

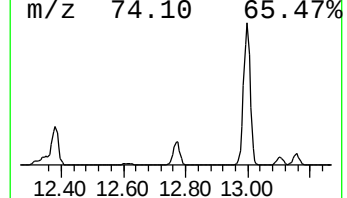
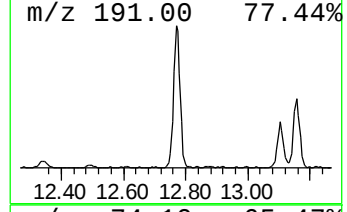
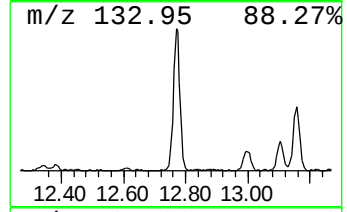
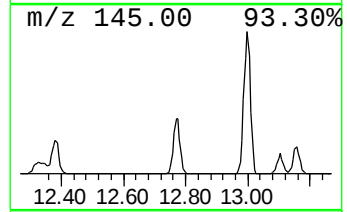
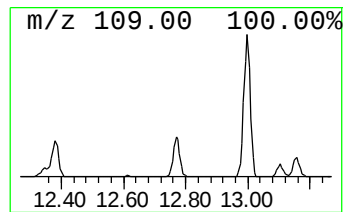
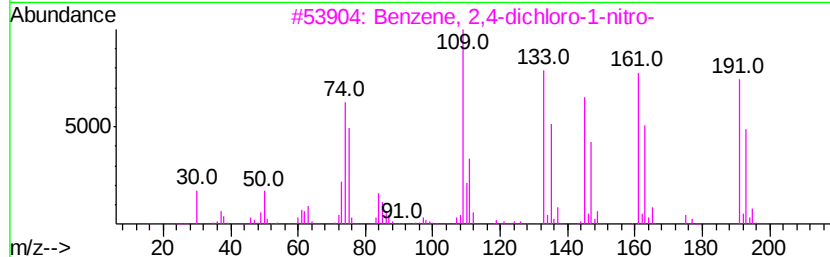
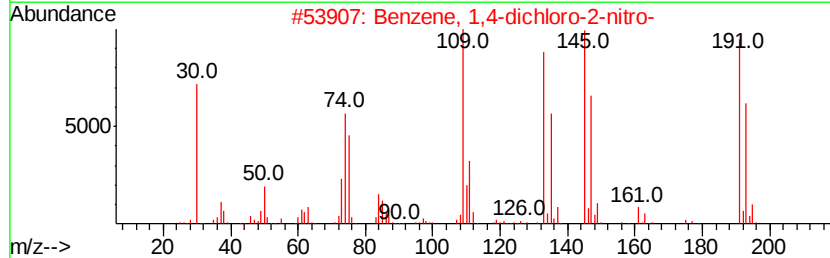
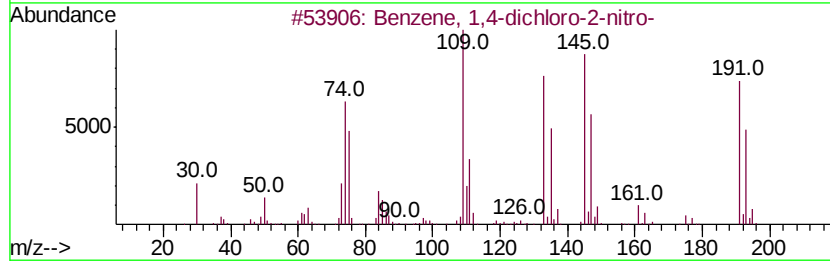
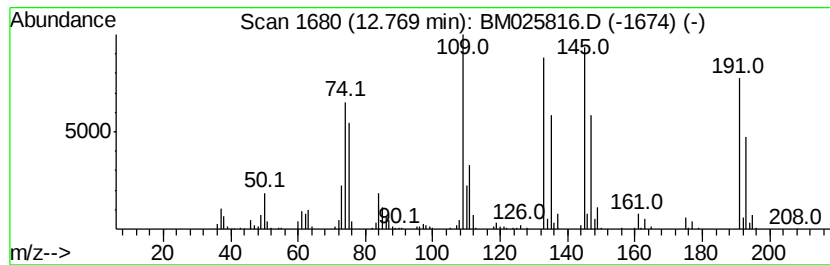
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1,4-dichloro-2-nitro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	7.39 ng/ul	856858	Acenaphthene-d10	14.20

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	98
3			Benzene, 2,4-dichloro-1-nitro-	191	C6H3Cl2NO2	000611-06-3	95
4			Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	95
5			Benzene, 2,4-dichloro-1-nitro-	191	C6H3Cl2NO2	000611-06-3	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025816.D
Acq On : 27 Mar 2020 00:55
Operator : CG/JU
Sample : L1968-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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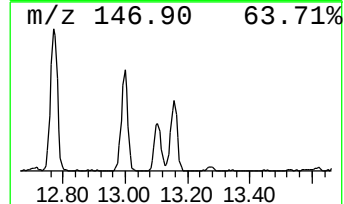
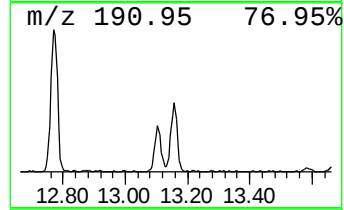
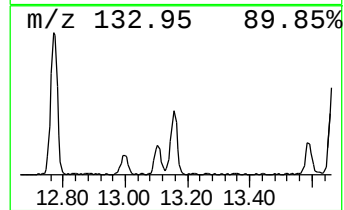
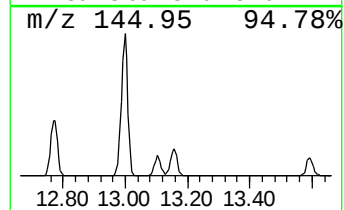
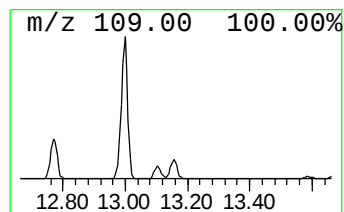
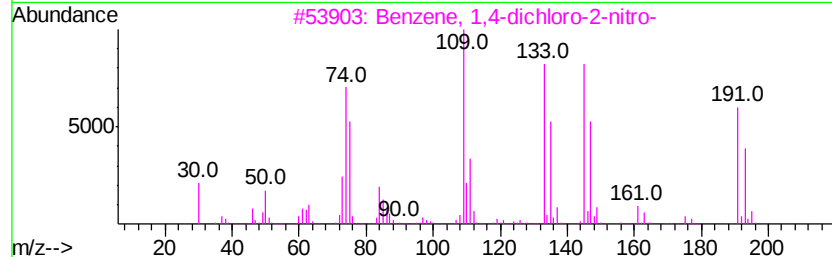
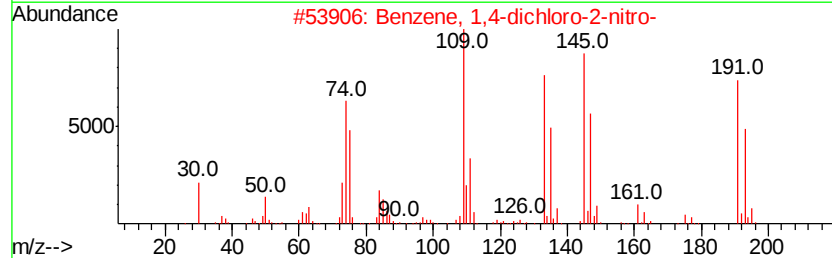
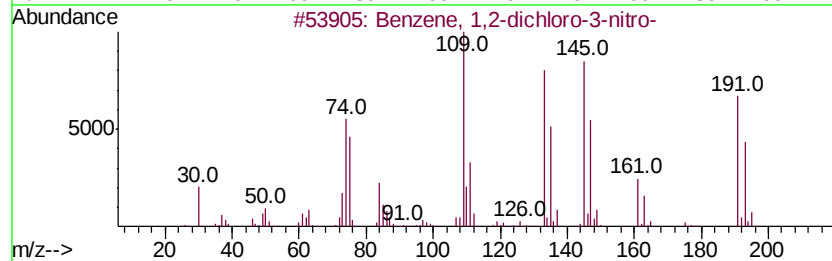
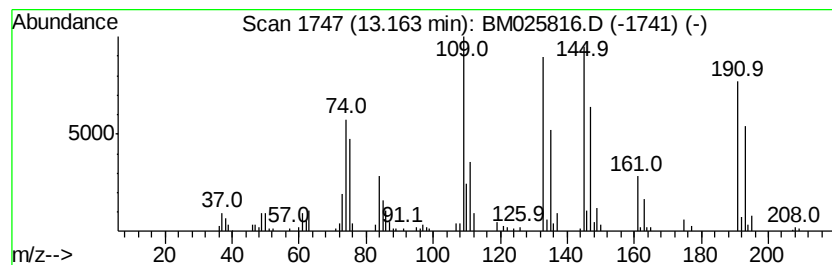
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 1,2-dichloro-3-nitro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	3.53 ng/ul	409147	Acenaphthene-d10	14.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-nitro-	191	C6H3Cl2NO2	003209-22-1	99
2		Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	99
3		Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	98
4		Benzene, 1,2-dichloro-3-nitro-	191	C6H3Cl2NO2	003209-22-1	97
5		Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025816.D
Acq On : 27 Mar 2020 00:55
Operator : CG/JU
Sample : L1968-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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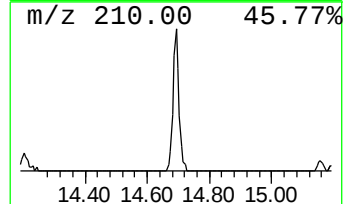
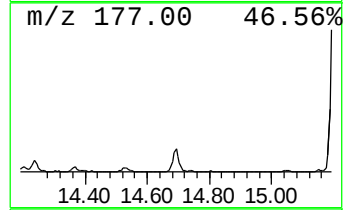
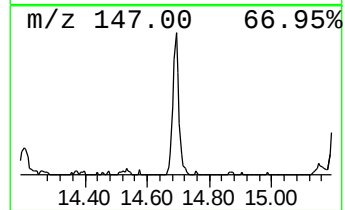
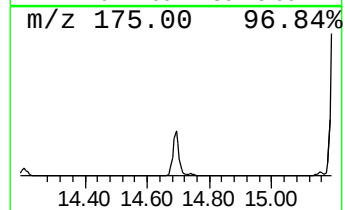
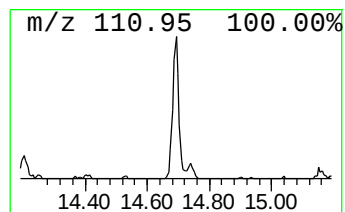
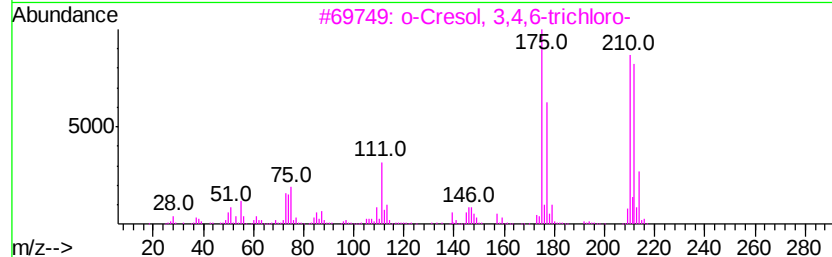
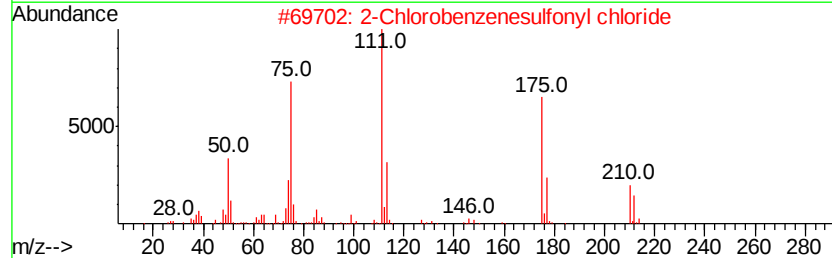
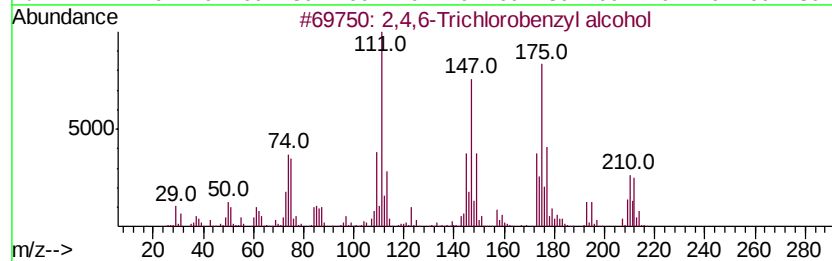
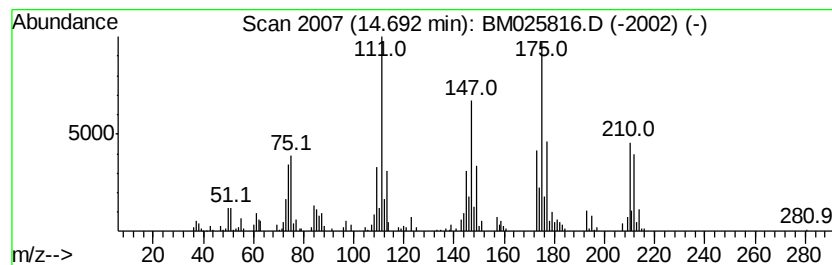
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 2,4,6-Trichlorobenzyl alcohol Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.69	3.35 ng/ul	388466	Acenaphthene-d10	14.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4,6-Trichlorobenzyl alcohol	210	C7H5Cl3O	217479-60-2	94
2			2-Chlorobenzenesulfonyl chloride	210	C6H4Cl2O2S	002905-23-9	64
3			o-Cresol, 3,4,6-trichloro-	210	C7H5Cl3O	000643-14-1	64
4			Benzenesulfonyl chloride, 4-chloro-	210	C6H4Cl2O2S	000098-60-2	58
5			Benzenemethanol, 2,3-dichloro-.a...	190	C8H8Cl2O	054798-91-3	50



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025816.D
Acq On : 27 Mar 2020 00:55
Operator : CG/JU
Sample : L1968-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

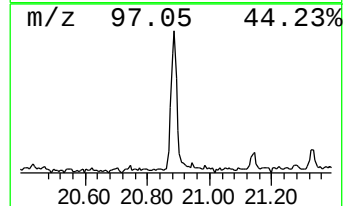
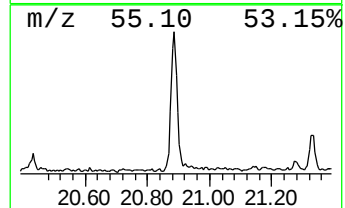
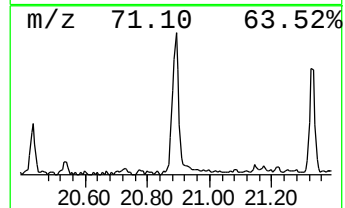
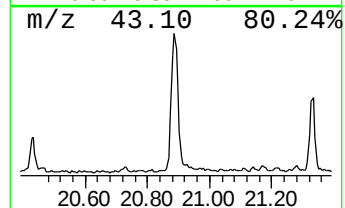
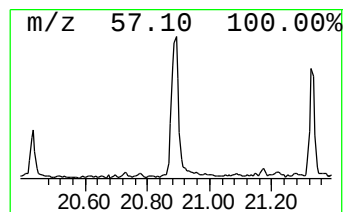
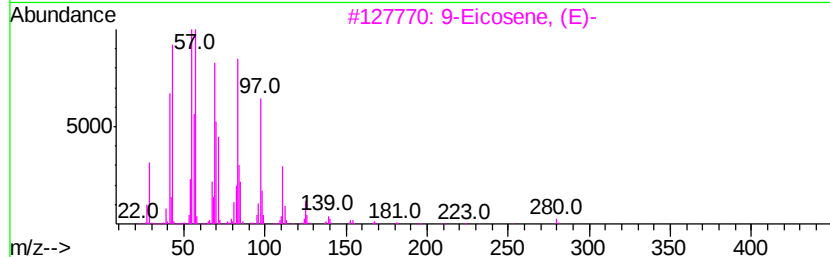
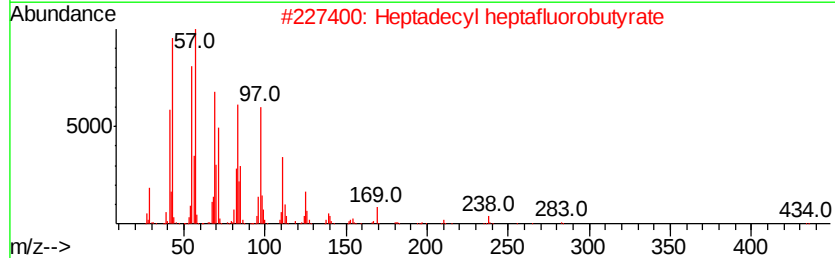
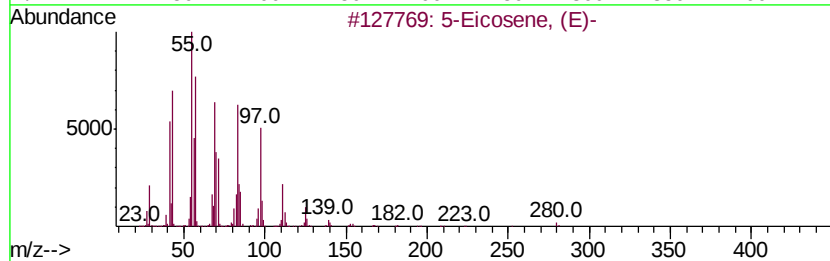
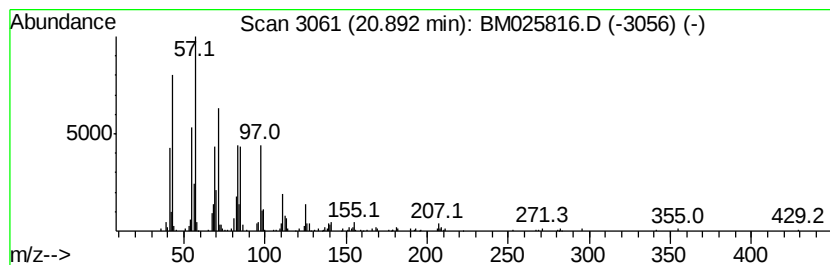
Instrument :
BNA_M
ClientSampleId :
CODE3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.		
20.89	3.00 ng/ul	481946	Chrysene-d12	21.14		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	97
2	Heptadecyl heptafluorobutyrte		452	C21H35F7O2	959085-66-6	94
3	9-Eicosene, (E)-		280	C20H40	074685-29-3	93
4	Trichloroacetic acid, hexadecyl ...		386	C18H33Cl3O2	074339-54-1	91
5	Pentafluoropropionic acid, hepta...		402	C20H35F5O2	959218-78-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032620\
Data File : BM025816.D
Acq On : 27 Mar 2020 00:55
Operator : CG/JU
Sample : L1968-12
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
CODE3

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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-bromo-...	9.05	11.6	ng/ul	946564	2	10.35	1637490	20.0
Benzene, 1-bromo-...	9.36	9.6	ng/ul	785816	2	10.35	1637490	20.0
Benzene, 1-chloro...	10.95	195.1	ng/ul	15976200	2	10.35	1637490	20.0
Benzene, 1-chloro...	11.15	51.6	ng/ul	4225950	2	10.35	1637490	20.0
4-Chloro-2-nitrop...	11.63	7.2	ng/ul	591280	2	10.35	1637490	20.0
Phenol, 5-chloro-...	12.32	8.4	ng/ul	980622	3	14.20	2319980	20.0
Benzene, 1,4-dich...	12.77	7.4	ng/ul	856858	3	14.20	2319980	20.0
Benzene, 1,2-dich...	13.16	3.5	ng/ul	409147	3	14.20	2319980	20.0
2,4,6-Trichlorobe...	14.69	3.4	ng/ul	388466	3	14.20	2319980	20.0
(DEL) Alkane: Str...	20.89	3.0	ng/ul	481946	5	21.14	3209080	20.0