

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
 Data File : BM016812.D
 Acq On : 20 Sep 2018 00:49
 Operator : SJ/JU
 Sample : J4896-16
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C08N7

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-BM091018MA.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	3.016	3	5	12	rVB	2524813	2877895	1.37%	0.277%
2	5.098	348	359	362	rBV3	51607430	137987271	65.54%	13.289%
3	6.422	578	584	593	rVB	292565	444621	0.21%	0.043%
4	6.898	658	665	672	rBV	1065351	1641397	0.78%	0.158%
5	7.075	689	695	705	rBV	727007	1283479	0.61%	0.124%
6	7.263	719	727	736	rBV	1043051	1740451	0.83%	0.168%
7	7.639	778	791	801	rBV2	43306086	108148627	51.37%	10.415%
8	7.804	801	819	821	rBV3	53816568	175699681	83.45%	16.921%
9	8.128	857	874	876	rBV4	53927385	181810837	86.36%	17.509%
10	8.428	918	925	933	rVB	1328806	2051789	0.97%	0.198%
11	8.881	995	1002	1005	rBV	961234	1592925	0.76%	0.153%
12	8.922	1005	1009	1017	rVB	870276	1362549	0.65%	0.131%
13	9.210	1051	1058	1070	rBV	3686688	6052341	2.87%	0.583%
14	9.439	1086	1097	1101	rBV	976332	1641114	0.78%	0.158%
15	9.480	1101	1104	1107	rVV	887180	1538165	0.73%	0.148%
16	9.522	1107	1111	1118	rVV	1979681	3465071	1.65%	0.334%
17	9.598	1118	1124	1129	rVV	708759	1236777	0.59%	0.119%
18	9.657	1129	1134	1143	rVB	441384	745712	0.35%	0.072%
19	9.892	1166	1174	1184	rVB	291559	542915	0.26%	0.052%
20	10.133	1205	1215	1222	rBV	1128402	2098633	1.00%	0.202%
21	10.422	1247	1264	1266	rBV6	58257148	210538278	100.00%	20.276%
22	10.533	1277	1283	1288	rBV	1754339	2558267	1.22%	0.246%
23	10.575	1288	1290	1296	rVB	174047	224561	0.11%	0.022%
24	10.651	1296	1303	1313	rVB	1050260	1865946	0.89%	0.180%
25	10.916	1336	1348	1354	rBV	46282214	102067324	48.48%	9.830%
26	11.104	1374	1380	1393	rVB	175186	297782	0.14%	0.029%
27	11.704	1473	1482	1488	rBV	212330	365652	0.17%	0.035%
28	11.792	1488	1497	1511	rVB4	251302	591743	0.28%	0.057%
29	12.074	1541	1545	1553	rVB2	142172	256347	0.12%	0.025%
30	12.251	1569	1575	1591	rBV2	168047	332155	0.16%	0.032%
31	12.504	1609	1618	1620	rBV	1149312	1833997	0.87%	0.177%
32	12.539	1620	1624	1631	rVV	4148742	6464447	3.07%	0.623%
33	12.774	1656	1664	1677	rVB	421384	685285	0.33%	0.066%
34	13.157	1721	1729	1741	rVB	16872849	25876258	12.29%	2.492%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Sampling : 1
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Max Peaks: 100
Peak Location: TOP

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Peak separation: 5

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35	13.427	1770	1775	1782	rVB2	164101	250522	0.12%	0.024%
36	13.739	1822	1828	1830	rBV2	257695	376057	0.18%	0.036%
37	13.774	1830	1834	1838	rVV	2099043	2886848	1.37%	0.278%
38	13.815	1838	1841	1854	rVB	801262	1188386	0.56%	0.114%
39	14.051	1872	1881	1889	rBV	2504011	3624529	1.72%	0.349%
40	14.363	1927	1934	1950	rVB2	2197678	3493068	1.66%	0.336%
41	14.510	1950	1959	1962	rBV5	147951	256203	0.12%	0.025%
42	14.557	1962	1967	1977	rVV	1023186	1467369	0.70%	0.141%
43	14.674	1981	1987	1996	rVB	562232	807992	0.38%	0.078%
44	14.845	2010	2016	2023	rBV	473684	719109	0.34%	0.069%
45	15.357	2096	2103	2116	rVB	3140542	4602789	2.19%	0.443%
46	15.468	2116	2122	2135	rBV	812157	1170283	0.56%	0.113%
47	17.109	2394	2401	2406	rBV	2645619	3728677	1.77%	0.359%
48	17.209	2411	2418	2434	rVV	3316010	4739515	2.25%	0.456%
49	19.503	2802	2808	2816	rBV	4404619	5634908	2.68%	0.543%
50	21.297	3107	3113	3121	rBV	3727368	4659567	2.21%	0.449%
51	23.391	3462	3469	3476	rBV	3239509	6038209	2.87%	0.582%
52	23.532	3486	3493	3509	rVB2	2517473	4808365	2.28%	0.463%

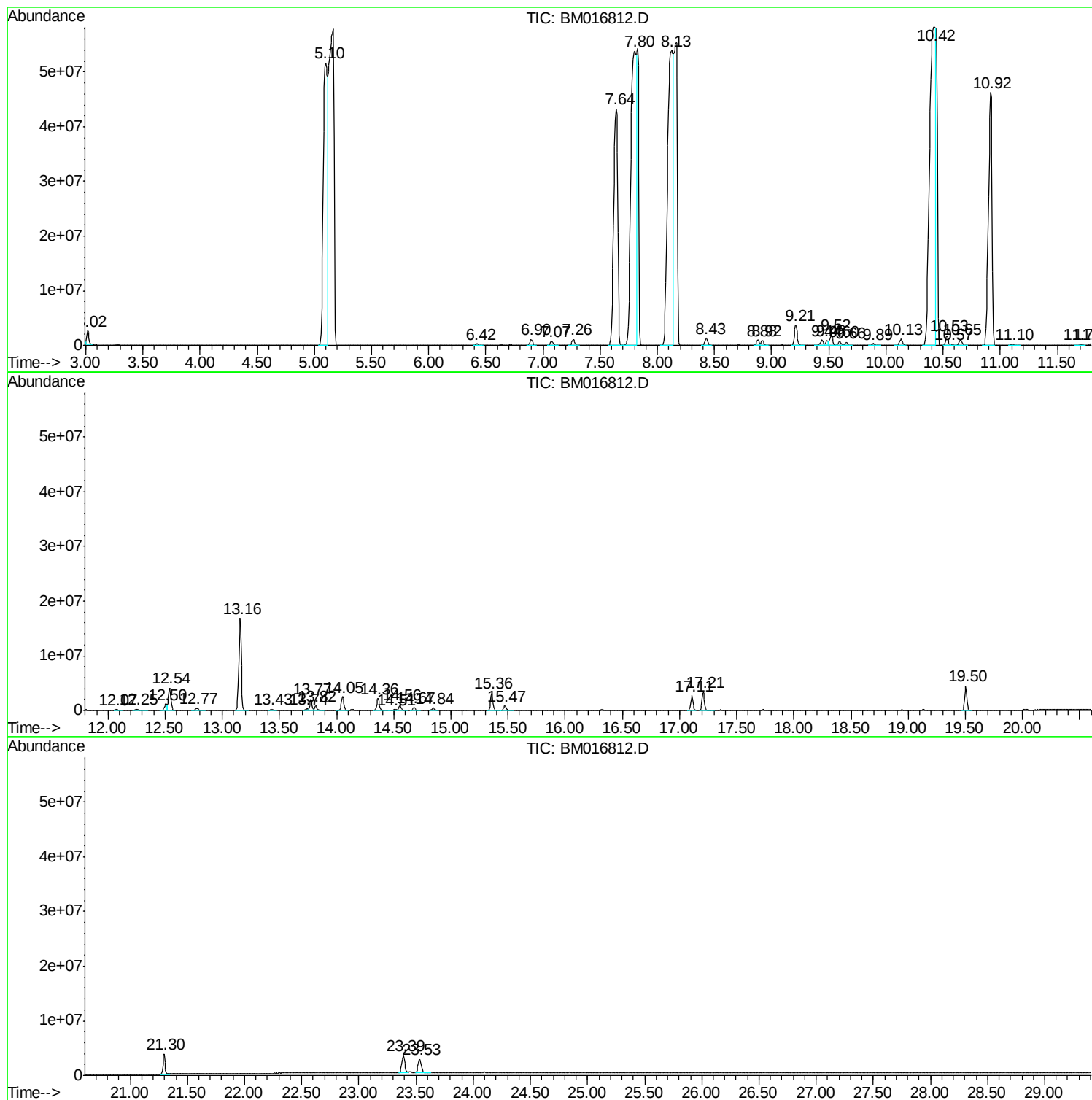
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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



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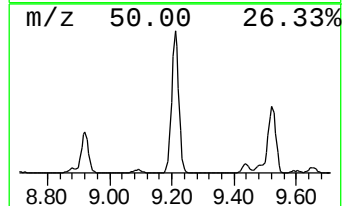
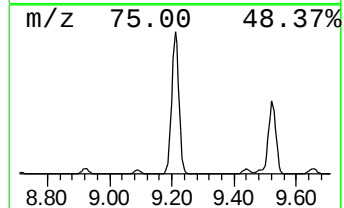
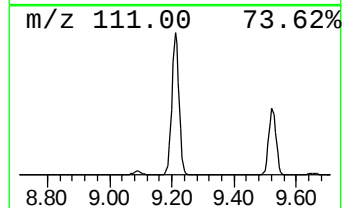
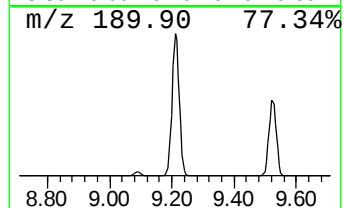
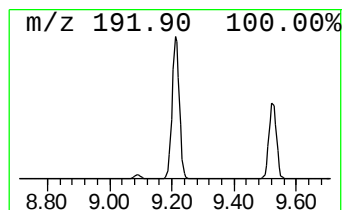
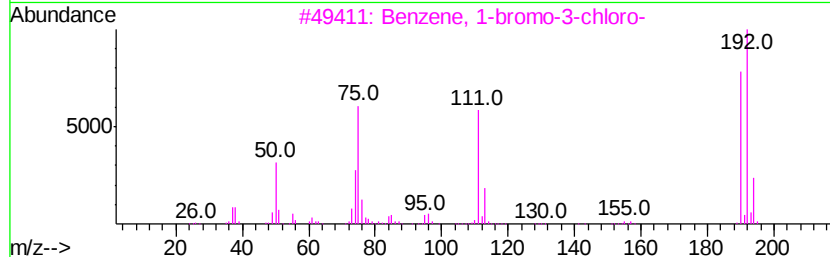
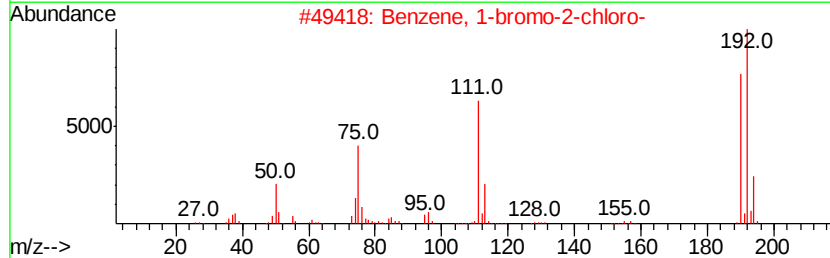
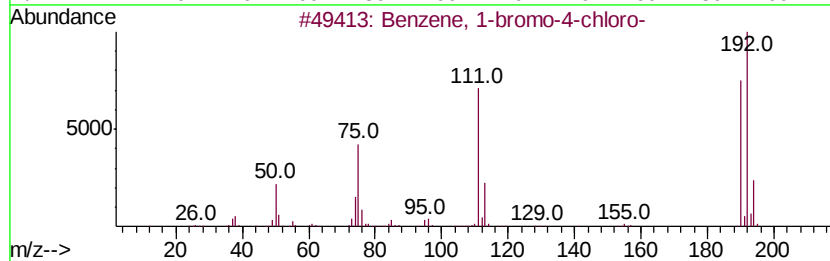
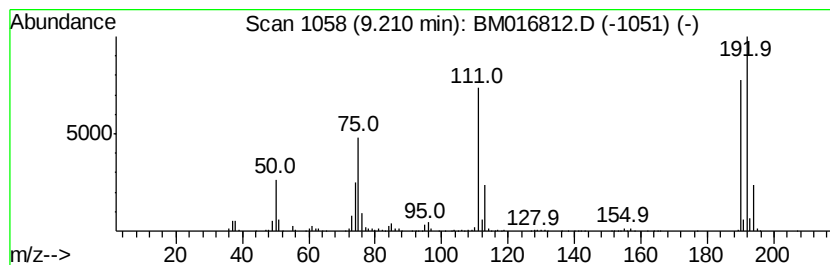
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1-bromo-4-chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.21	47.32 ng/ul	6052340	Napthalene-d8	10.53

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



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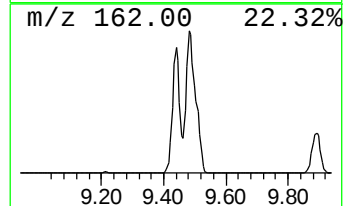
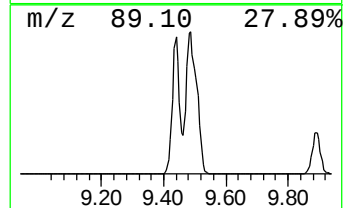
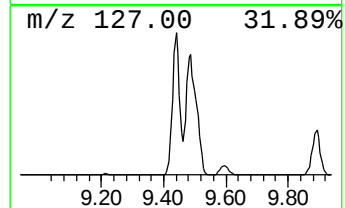
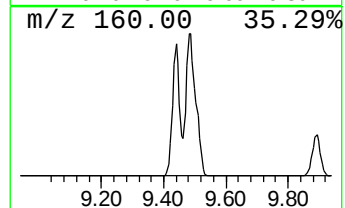
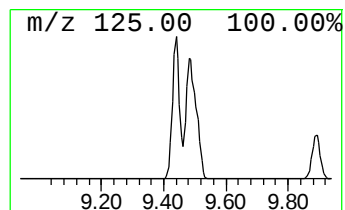
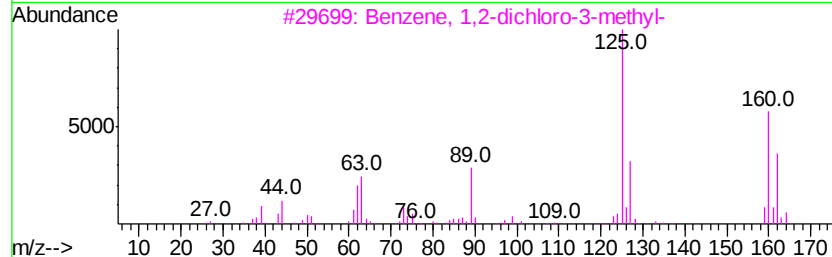
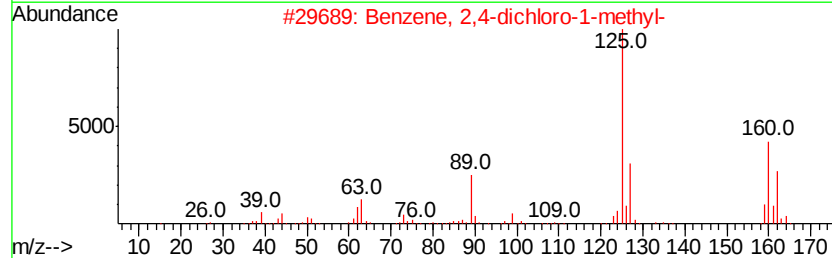
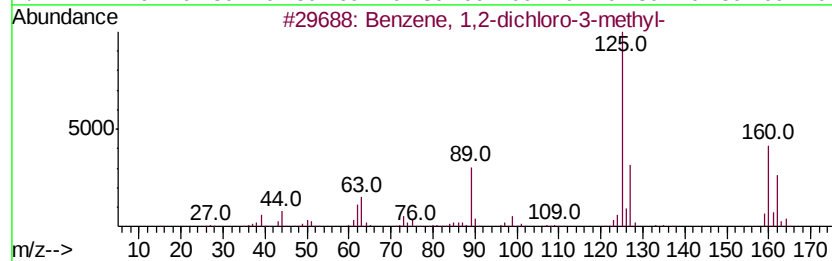
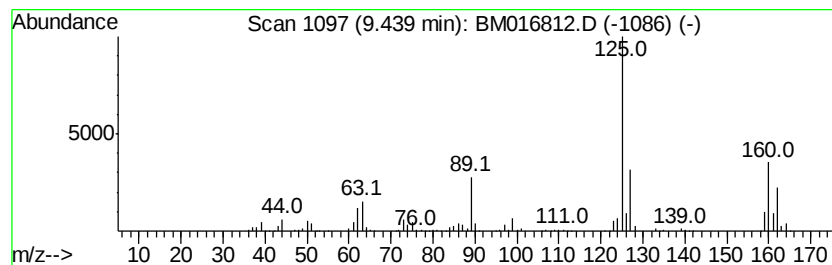
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1,2-dichloro-3-met... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	12.83 ng/ul	1641110	Napthalene-d8	10.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2		Benzene, 2,4-dichloro-1-methyl-	160	C7H6Cl2	000095-73-8	98
3		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
4		Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
5		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97



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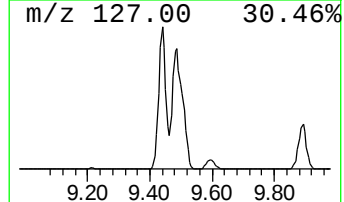
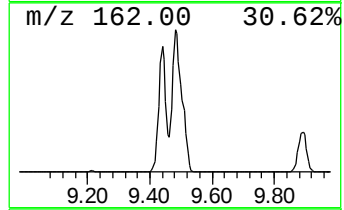
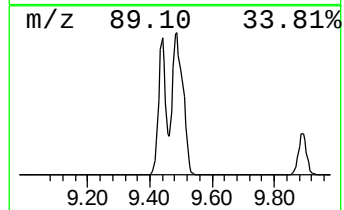
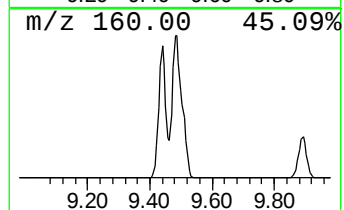
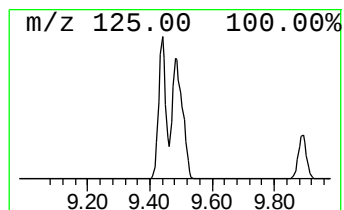
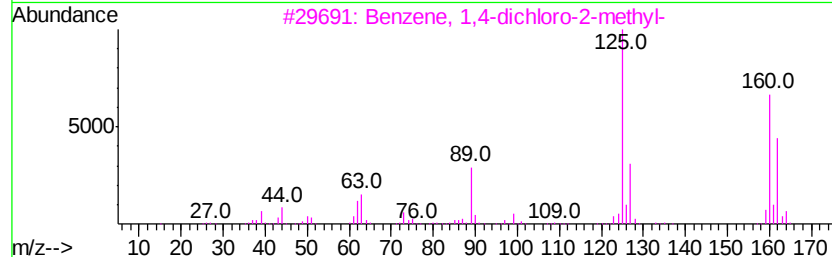
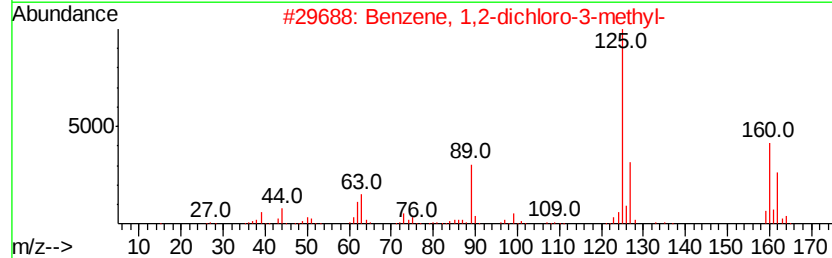
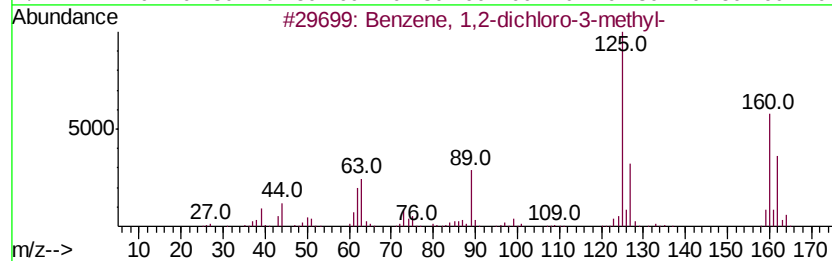
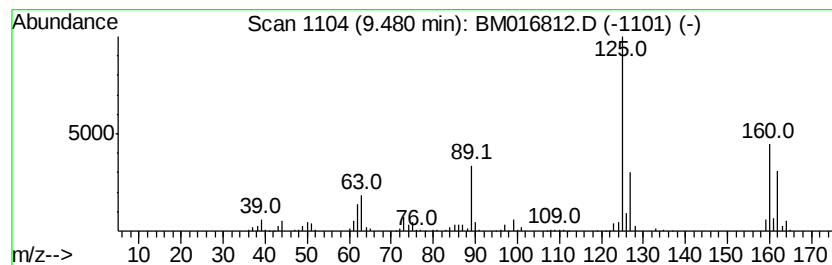
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,4-dichloro-2-met... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	12.03 ng/ul	1538170	Napthalene-d8	10.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
2		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96
3		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
4		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
5		Benzene, 1-chloro-2-(chloromethyl)-	160	C7H6Cl2	000611-19-8	96



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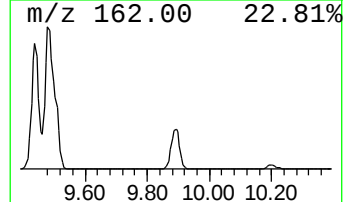
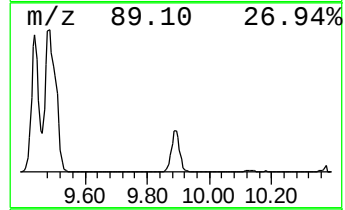
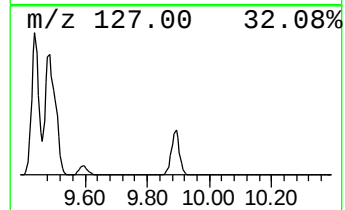
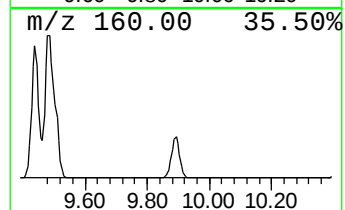
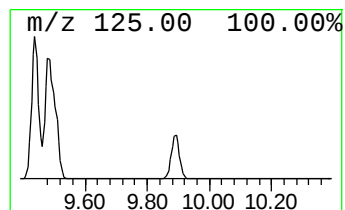
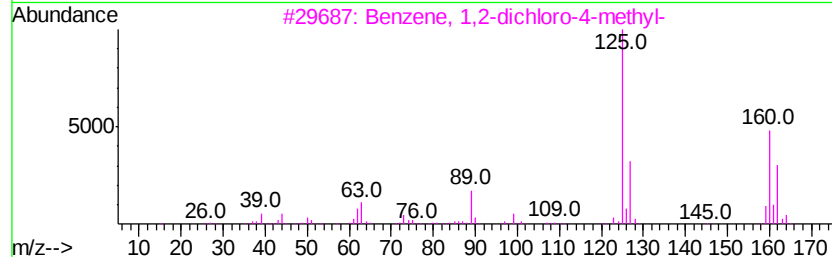
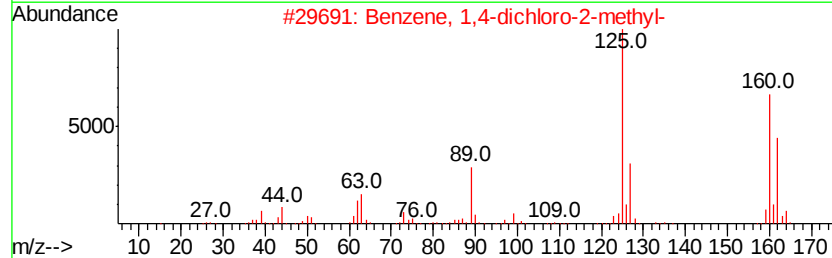
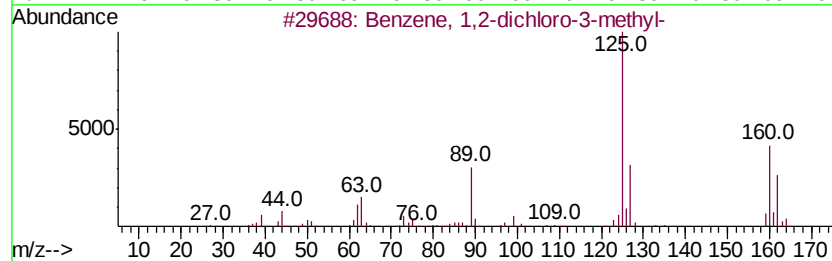
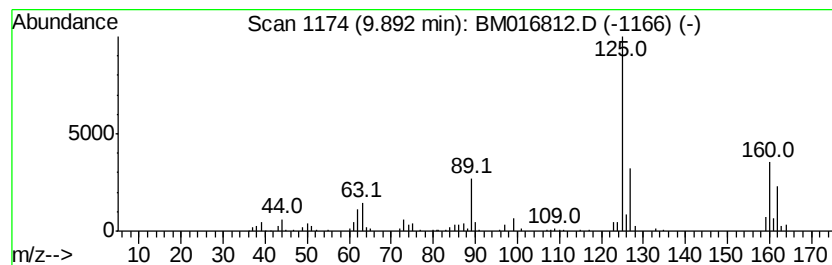
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 8 Benzene, 1,2-dichloro-4-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.89	4.24 ng/ul	542915	Naphthalene-d8	10.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	97
3		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97
4		Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	96
5		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96



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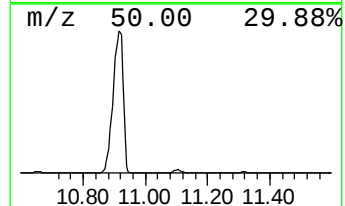
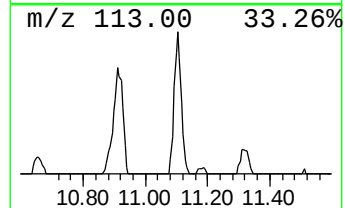
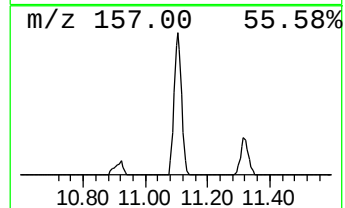
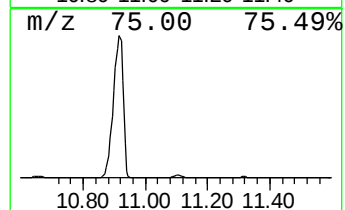
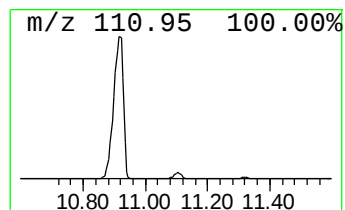
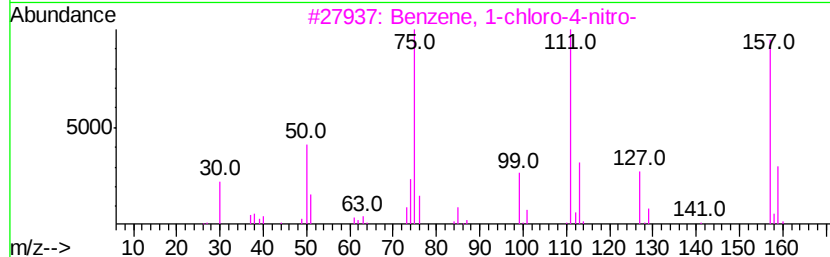
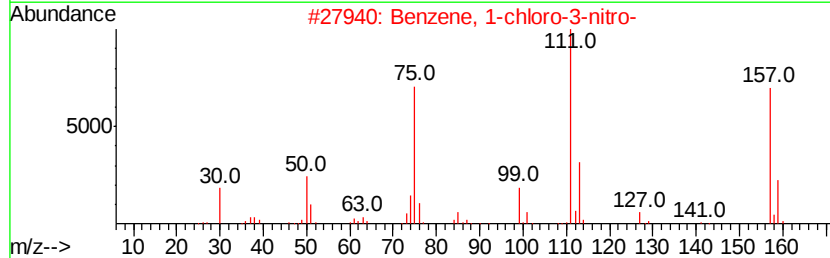
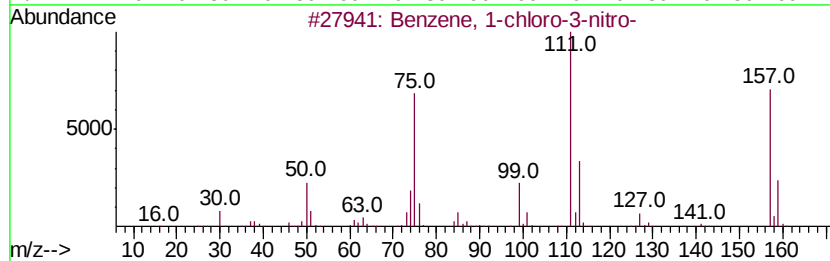
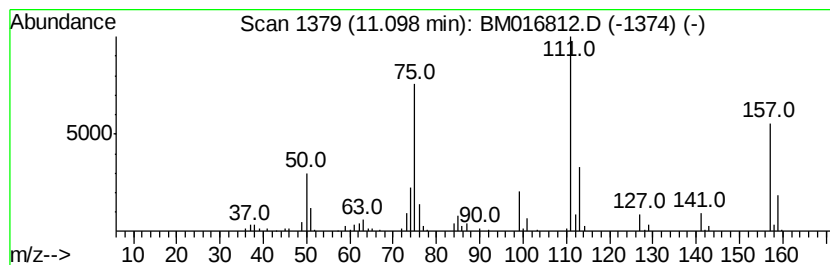
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1-chloro-3-nitro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.10	2.33 ng/ul	297782	Naphthalene-d8	10.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
2			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
3			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
4			Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	92
5			Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
Operator : SJ/JU
Sample : J4896-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

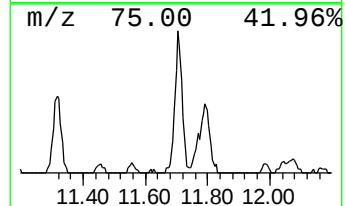
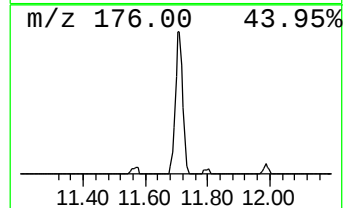
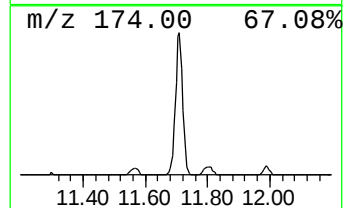
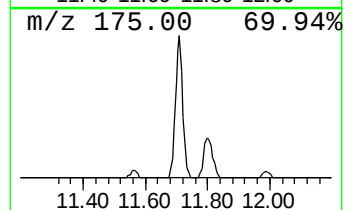
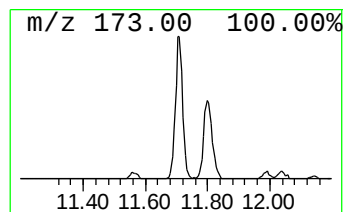
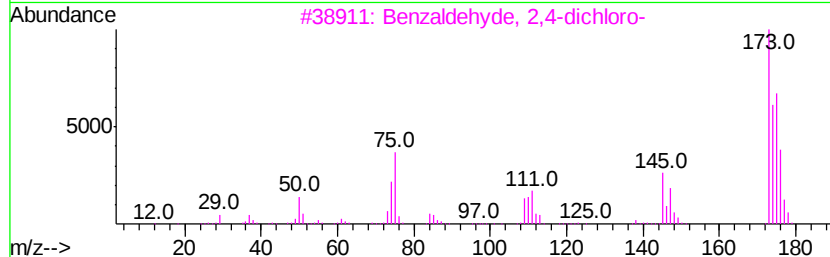
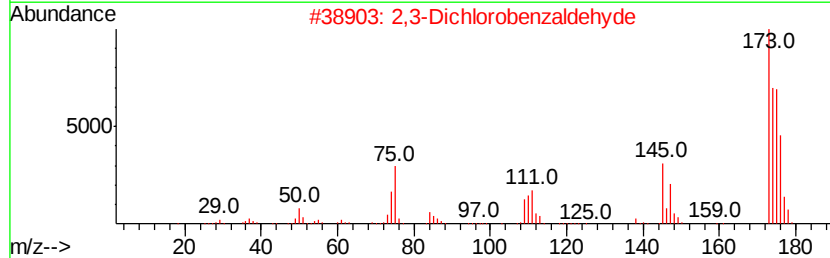
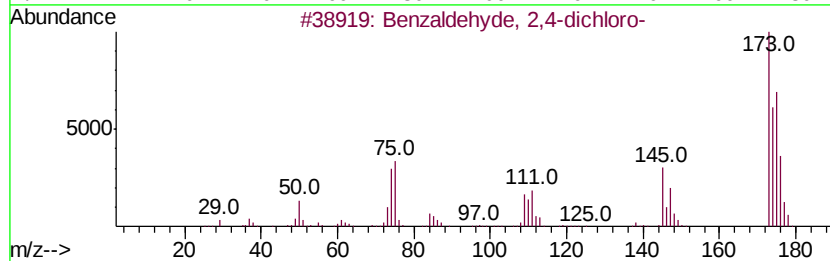
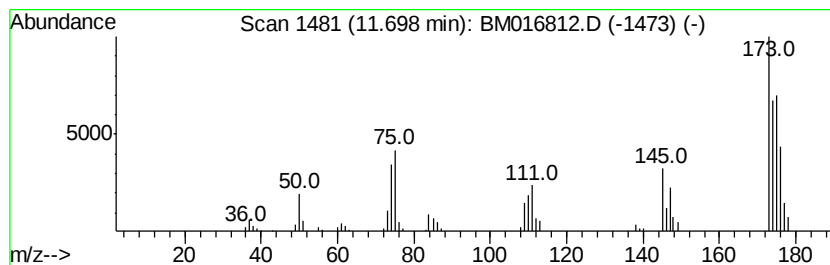
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzaldehyde, 2,4-dichloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	2.86 ng/ul	365652	Napthalene-d8	10.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzaldehyde, 2,4-dichloro-	174 C7H4Cl2O	000874-42-0 99
2		2,3-Dichlorobenzaldehyde	174 C7H4Cl2O	006334-18-5 99
3		Benzaldehyde, 2,4-dichloro-	174 C7H4Cl2O	000874-42-0 99
4		2,3-Dichlorobenzaldehyde	174 C7H4Cl2O	006334-18-5 99
5		3,5-Dichlorobenzaldehyde	174 C7H4Cl2O	010203-08-4 98



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
Operator : SJ/JU
Sample : J4896-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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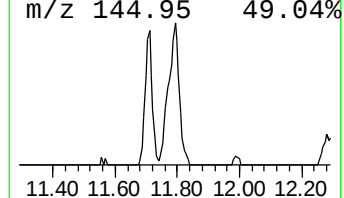
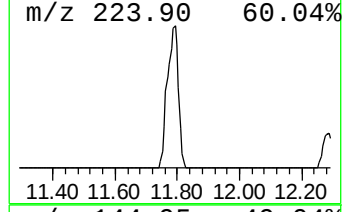
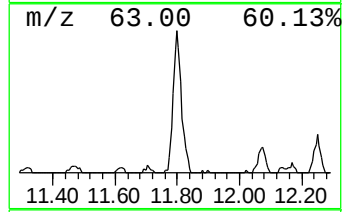
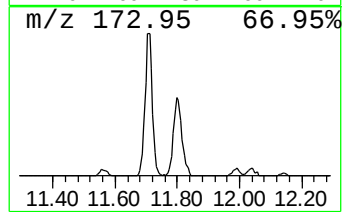
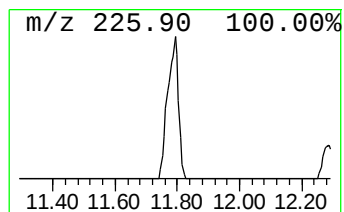
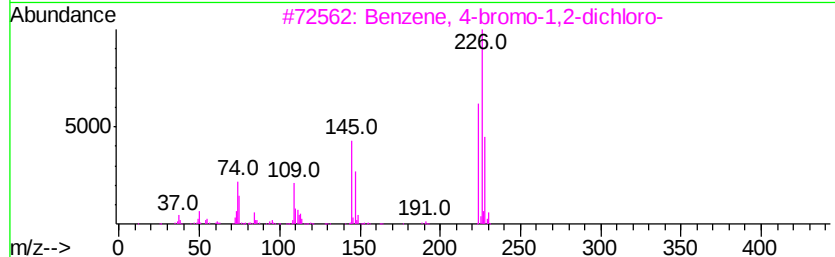
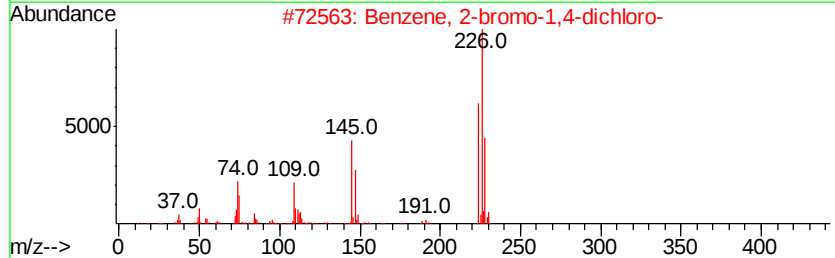
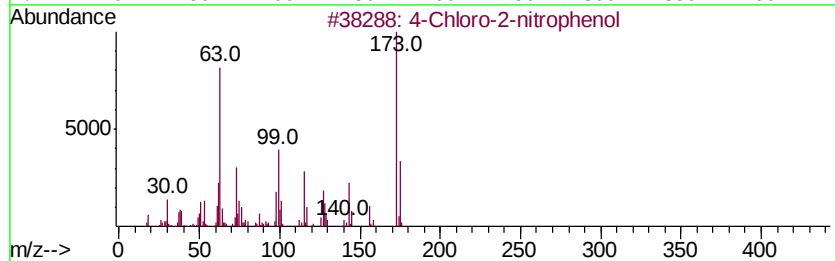
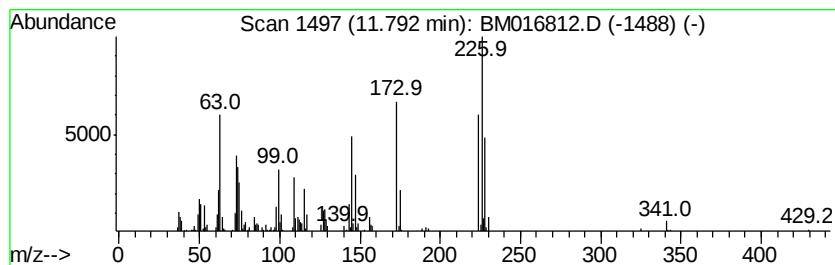
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 4-Chloro-2-nitrophenol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.79	4.63 ng/ul	591743	Naphthalene-d8	10.53

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	95
2		Benzene, 2-bromo-1,4-dichloro-	224	C6H3BrCl2	001435-50-3	95
3		Benzene, 4-bromo-1,2-dichloro-	224	C6H3BrCl2	018282-59-2	95
4		1-Bromo-2,4-dichlorobenzene	224	C6H3BrCl2	001193-72-2	95
5		Benzene, 1-bromo-2,6-dichloro-	224	C6H3BrCl2	019393-92-1	92



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
Operator : SJ/JU
Sample : J4896-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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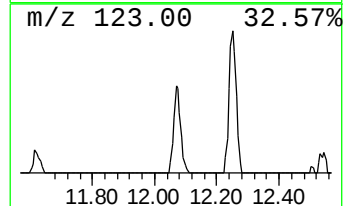
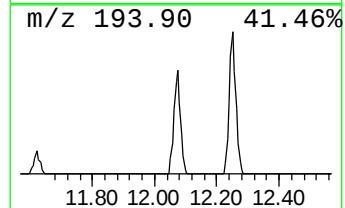
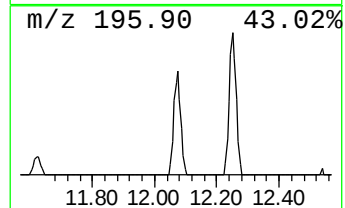
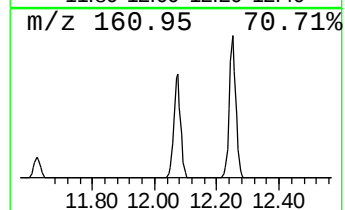
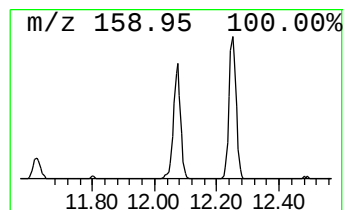
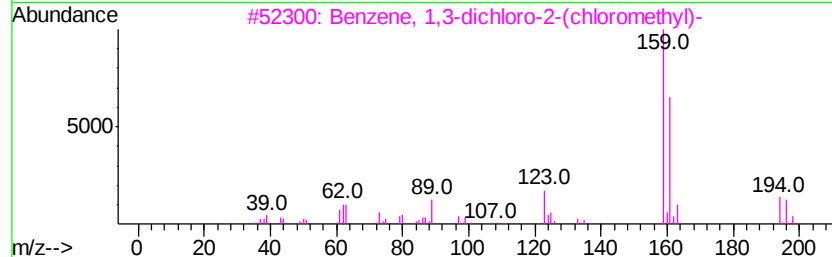
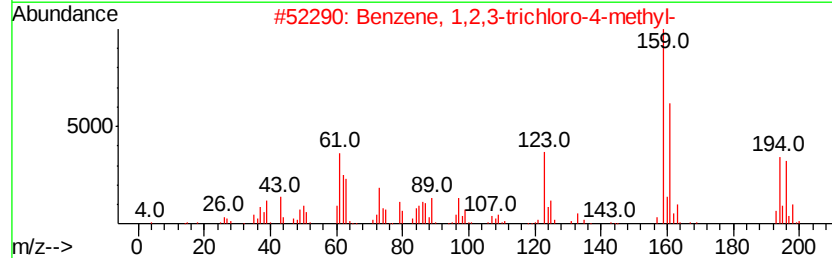
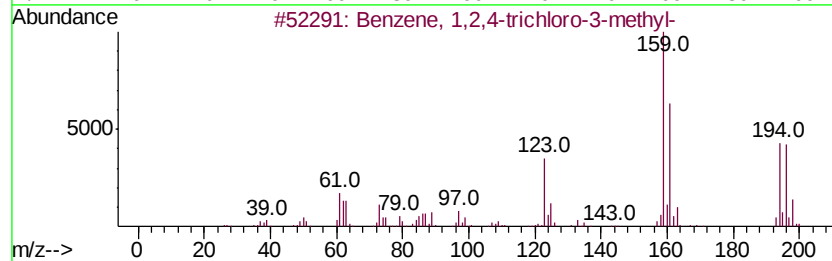
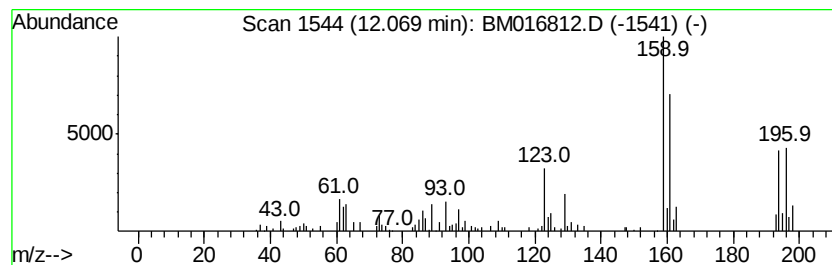
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1,2,4-trichloro-3-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.07	2.00 ng/ul	256347	Naphthalene-d8	10.53

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trichloro-3-methyl-	194	C7H5Cl3	002077-46-5	98
2			Benzene, 1,2,3-trichloro-4-methyl-	194	C7H5Cl3	007359-72-0	93
3			Benzene, 1,3-dichloro-2-(chloromethyl)-	194	C7H5Cl3	002014-83-7	62
4			Benzene, 1,2-dichloro-3-(chloromethyl)-	194	C7H5Cl3	003290-01-5	58
5			Benzene, 1,2-dichloro-4-(chloromethyl)-	194	C7H5Cl3	000102-47-6	58



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
Operator : SJ/JU
Sample : J4896-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

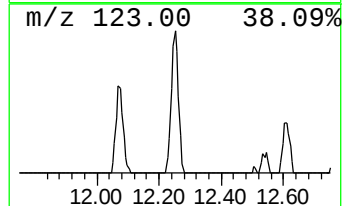
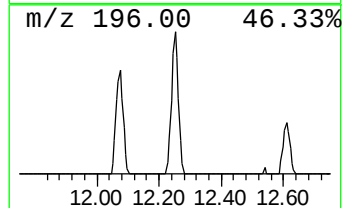
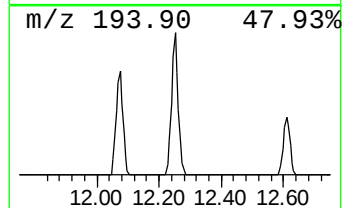
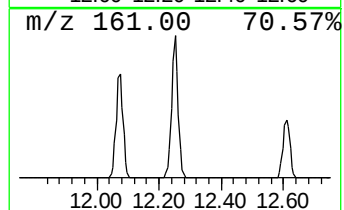
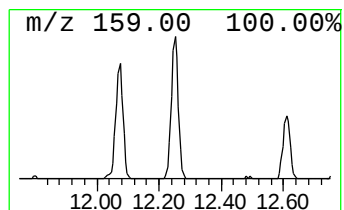
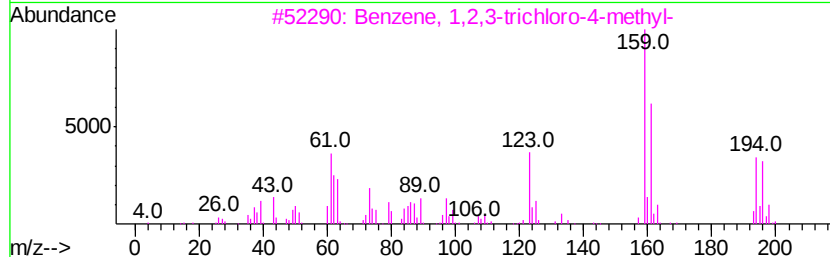
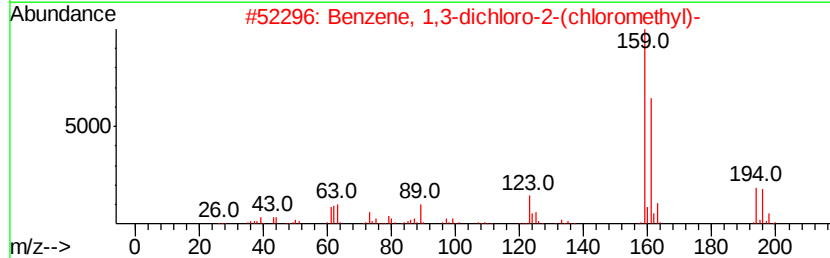
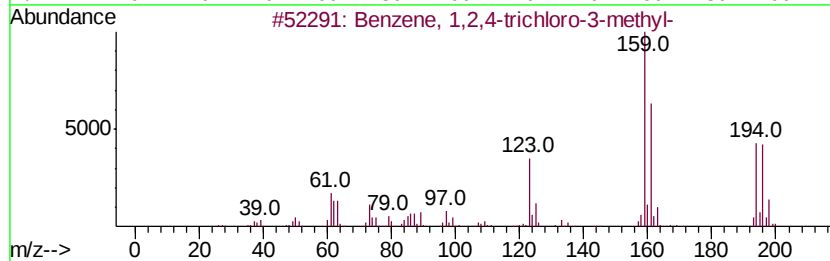
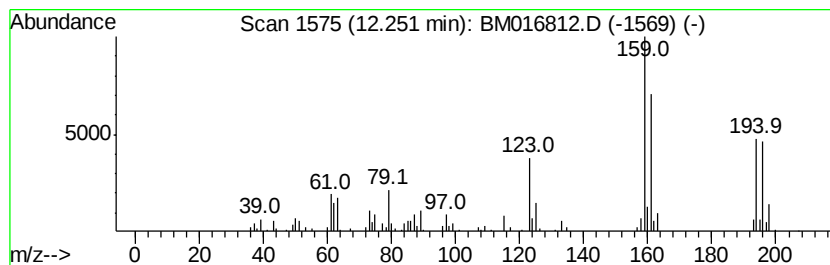
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1,3-dichloro-2-(ch... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	2.60 ng/ul	332155	Naphthalene-d8	10.53
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,4-trichloro-3-methyl-	194 C7H5Cl3	002077-46-5 99
2		Benzene, 1,3-dichloro-2-(chlorom...	194 C7H5Cl3	002014-83-7 96
3		Benzene, 1,2,3-trichloro-4-methyl-	194 C7H5Cl3	007359-72-0 95
4		2,4,5-Trichlorotoluene	194 C7H5Cl3	006639-30-1 95
5		Benzene, 1,2-dichloro-4-(chlorom...	194 C7H5Cl3	000102-47-6 93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
Operator : SJ/JU
Sample : J4896-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
C08N7

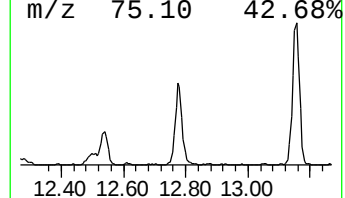
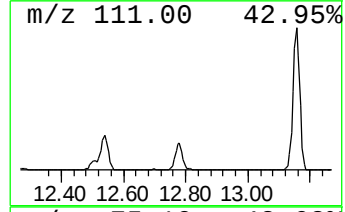
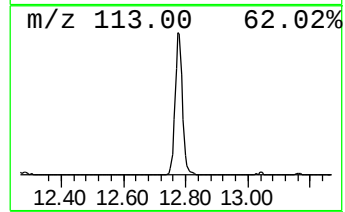
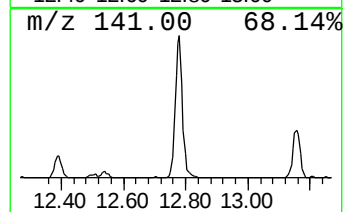
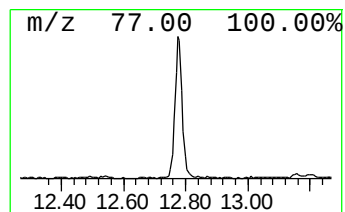
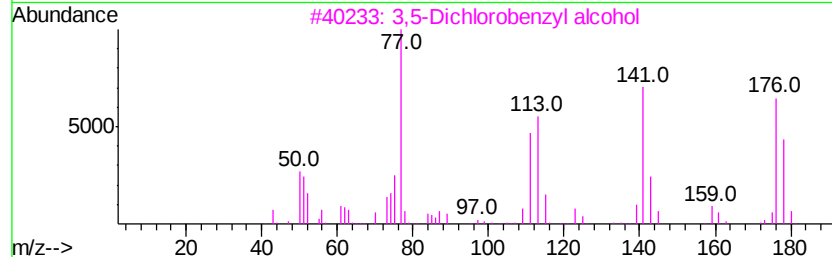
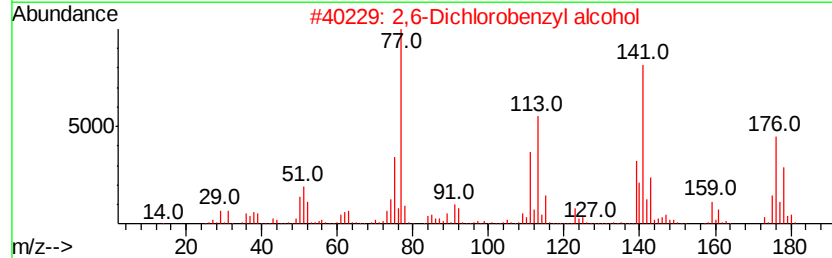
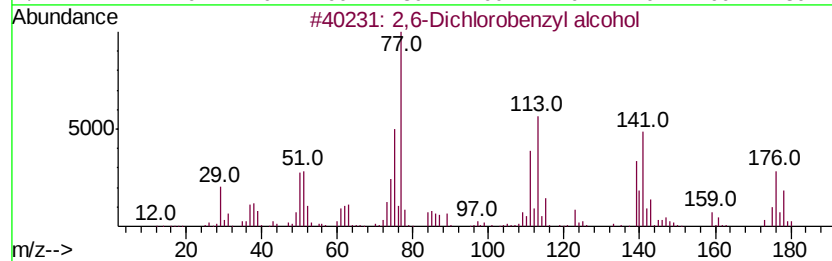
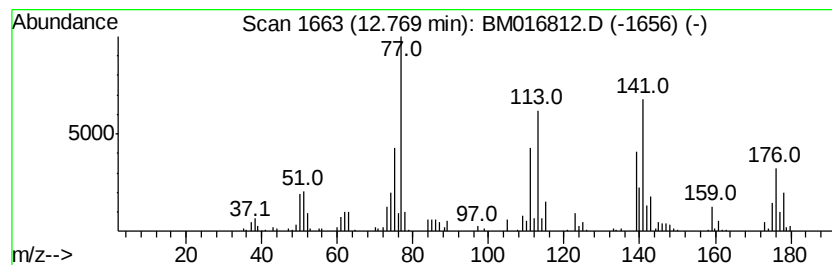
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 2,6-Dichlorobenzyl alcohol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	3.92 ng/ul	685285	Acenaphthene-d10	14.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,6-Dichlorobenzyl alcohol	176	C7H6Cl2O	015258-73-8	97
2			2,6-Dichlorobenzyl alcohol	176	C7H6Cl2O	015258-73-8	95
3			3,5-Dichlorobenzyl alcohol	176	C7H6Cl2O	060211-57-6	93
4			Benzenemethanol, 2,4-dichloro-	176	C7H6Cl2O	001777-82-8	93
5			3,5-Dichlorobenzyl alcohol	176	C7H6Cl2O	060211-57-6	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
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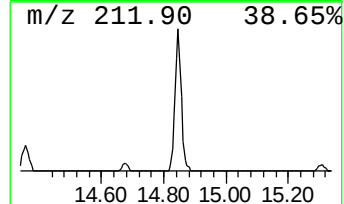
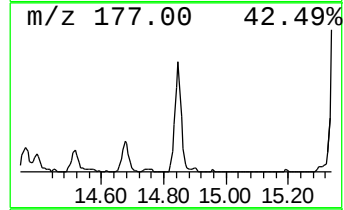
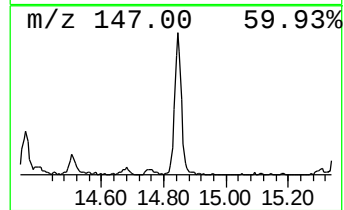
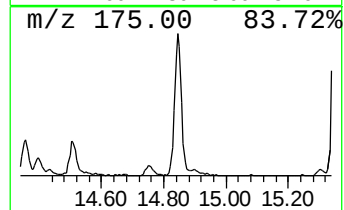
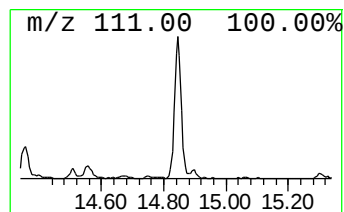
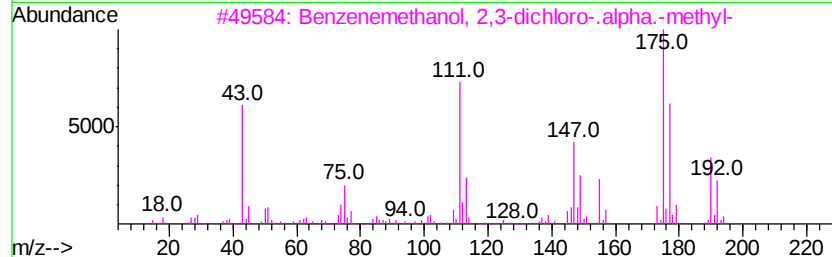
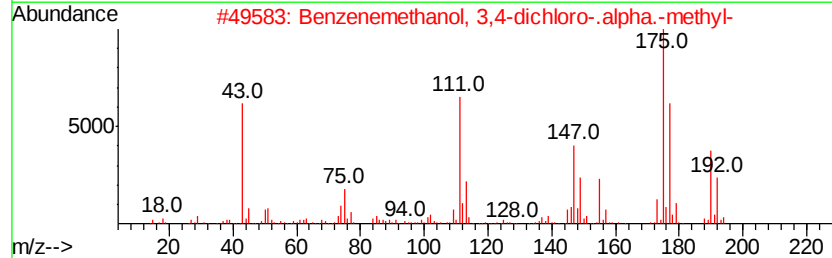
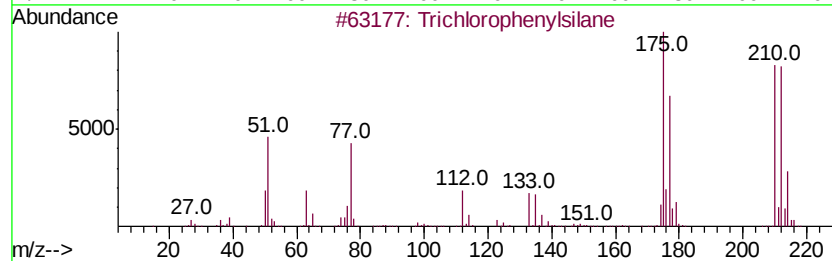
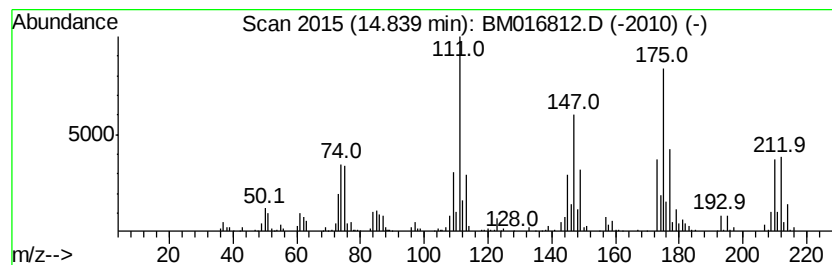
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Trichlorophenylsilane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	4.12 ng/ul	719109	Acenaphthene-d10	14.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Trichlorophenylsilane	210	C6H5Cl3Si	000098-13-5	50
2			Benzenemethanol, 3,4-dichloro-.a...	190	C8H8Cl2O	001475-11-2	50
3			Benzenemethanol, 2,3-dichloro-.a...	190	C8H8Cl2O	054798-91-3	50
4			2-Chlorobenzenesulfonyl chloride	210	C6H4Cl2O2S	002905-23-9	43
5			Benzenemethanol, 2,4-dichloro-.a...	224	C8H7Cl3O	013692-14-3	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM091918\
Data File : BM016812.D
Acq On : 20 Sep 2018 00:49
Operator : SJ/JU
Sample : J4896-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C08N7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.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, 1-bromo-...	9.21	47.3	ng/ul	6052340	2	10.53	2558270	20.0
Benzene, 1,2-dich...	9.44	12.8	ng/ul	1641110	2	10.53	2558270	20.0
Benzene, 1,4-dich...	9.48	12.0	ng/ul	1538170	2	10.53	2558270	20.0
Benzene, 1,2-dich...	9.89	4.2	ng/ul	542915	2	10.53	2558270	20.0
Benzene, 1-chloro...	11.10	2.3	ng/ul	297782	2	10.53	2558270	20.0
Benzaldehyde, 2,4...	11.70	2.9	ng/ul	365652	2	10.53	2558270	20.0
4-Chloro-2-nitrop...	11.79	4.6	ng/ul	591743	2	10.53	2558270	20.0
Benzene, 1,2,4-tr...	12.07	2.0	ng/ul	256347	2	10.53	2558270	20.0
Benzene, 1,3-dich...	12.25	2.6	ng/ul	332155	2	10.53	2558270	20.0
2,6-Dichlorobenzy...	12.77	3.9	ng/ul	685285	3	14.36	3493070	20.0
Trichlorophenylsi...	14.84	4.1	ng/ul	719109	3	14.36	3493070	20.0