

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025820.D
 Acq On : 27 Mar 2020 03:56
 Operator : CG/JU
 Sample : L1968-16
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0DL6

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	4.987	340	357	365	rBV7	68114849	406747585	100.00%	25.137%
2	6.269	569	575	583	rVB	333449	520649	0.13%	0.032%
3	6.934	678	688	699	rBV	406089	919596	0.23%	0.057%
4	7.116	710	719	732	rVB	573780	1113532	0.27%	0.069%
5	7.487	768	782	794	rBV	24039110	76943877	18.92%	4.755%
6	7.681	794	815	819	rBV7	73289058	380935289	93.65%	23.542%
7	8.028	851	874	875	rBV6	73257607	373826061	91.91%	23.102%
8	8.281	911	917	928	rVB	567288	881031	0.22%	0.054%
9	8.716	984	991	994	rBV	719586	1222654	0.30%	0.076%
10	8.769	994	1000	1010	rVB	6540975	10682647	2.63%	0.660%
11	9.045	1039	1047	1055	rBV	1263044	2045620	0.50%	0.126%
12	9.281	1081	1087	1090	rBV	301613	463503	0.11%	0.029%
13	9.322	1090	1094	1096	rVV	279485	435313	0.11%	0.027%
14	9.357	1096	1100	1107	rVV	707608	1205606	0.30%	0.075%
15	9.428	1107	1112	1120	rVV2	638193	1212967	0.30%	0.075%
16	9.969	1197	1204	1212	rBV	913682	1559725	0.38%	0.096%
17	10.263	1237	1254	1259	rBV3	68162442	219646347	54.00%	13.574%
18	10.357	1265	1270	1276	rBV	999586	1412525	0.35%	0.087%
19	10.745	1324	1336	1342	rBV	34632272	67700571	16.64%	4.184%
20	10.939	1360	1369	1384	rVB	4884299	7852328	1.93%	0.485%
21	11.145	1395	1404	1414	rVB	998497	1670165	0.41%	0.103%
22	11.627	1476	1486	1492	rBV3	225580	432317	0.11%	0.027%
23	12.321	1597	1604	1606	rBV	283125	462480	0.11%	0.029%
24	12.345	1606	1608	1610	rVV	378103	496393	0.12%	0.031%
25	12.380	1610	1614	1621	rVB	1944838	2886496	0.71%	0.178%
26	12.998	1712	1719	1726	rVB	7586169	11066340	2.72%	0.684%
27	13.592	1814	1820	1821	rBV	344745	432489	0.11%	0.027%
28	13.621	1821	1825	1830	rVV	1844810	2586218	0.64%	0.160%
29	13.892	1865	1871	1881	rVB	2174420	3082357	0.76%	0.190%
30	14.204	1915	1924	1933	rBV2	1348347	1941864	0.48%	0.120%
31	15.204	2087	2094	2100	rBV	2702640	3807522	0.94%	0.235%
32	15.327	2109	2115	2123	rBV	876463	1178386	0.29%	0.073%
33	16.951	2385	2391	2397	rBV	1682399	2259102	0.56%	0.140%
34	17.051	2401	2408	2419	rVB	3028519	4170450	1.03%	0.258%

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35	18.327	2620	2625	2631	rVV	5380692	7140331	1.76%	0.441%
36	18.656	2677	2681	2685	rVV	342544	487285	0.12%	0.030%
37	19.345	2792	2798	2805	rBV	3615391	4925968	1.21%	0.304%
38	20.886	3055	3060	3067	rBV	361858	493784	0.12%	0.031%
39	21.139	3098	3103	3110	rBV	2221615	2736737	0.67%	0.169%
40	23.156	3439	3446	3453	rBV	3330094	5615797	1.38%	0.347%
41	23.291	3462	3469	3477	rVB	1651914	2922121	0.72%	0.181%

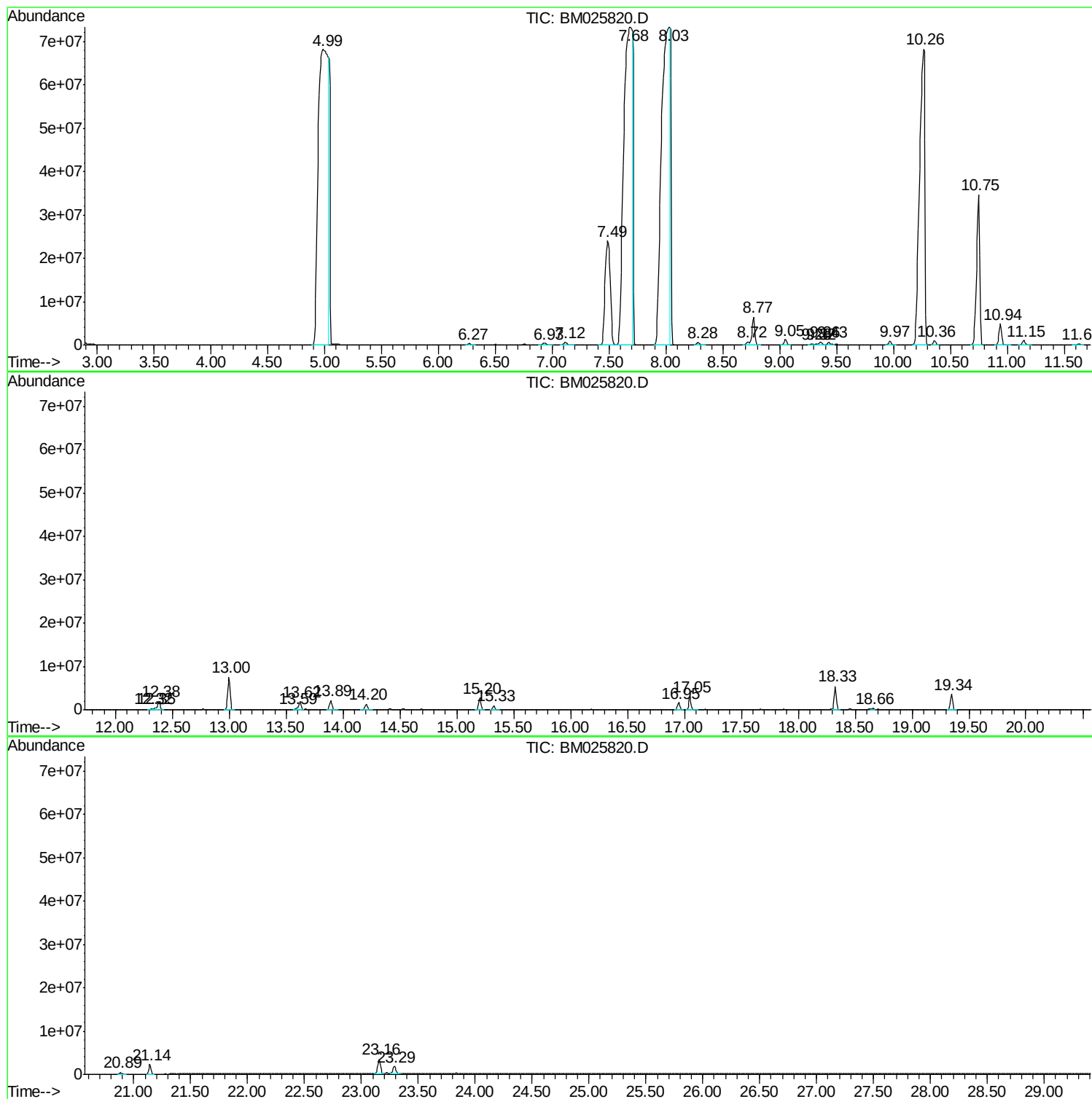
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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



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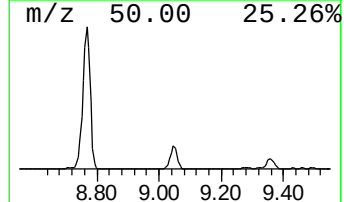
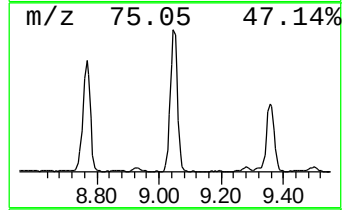
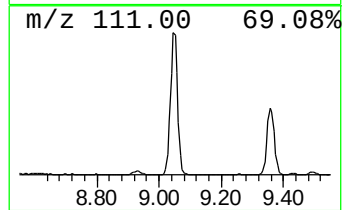
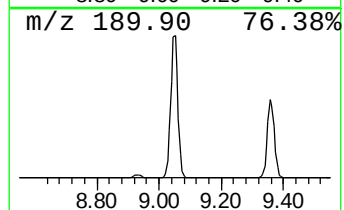
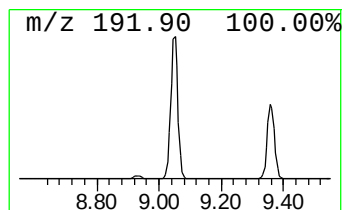
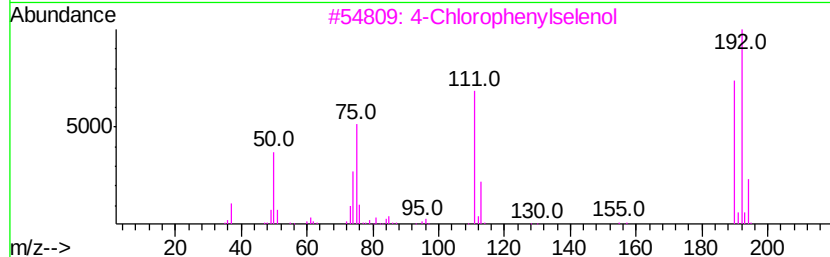
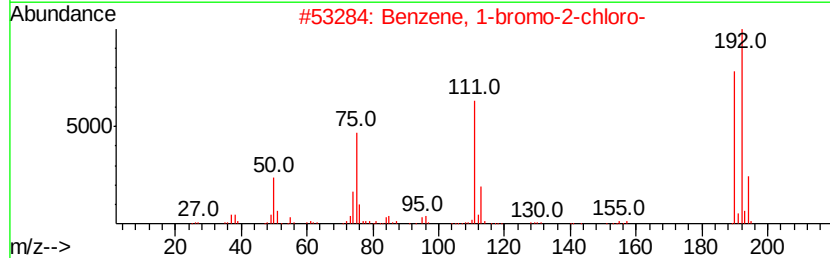
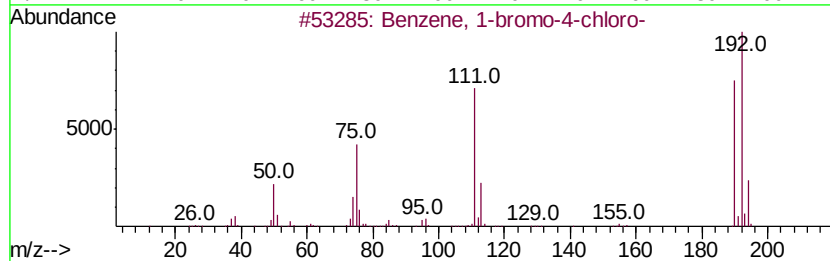
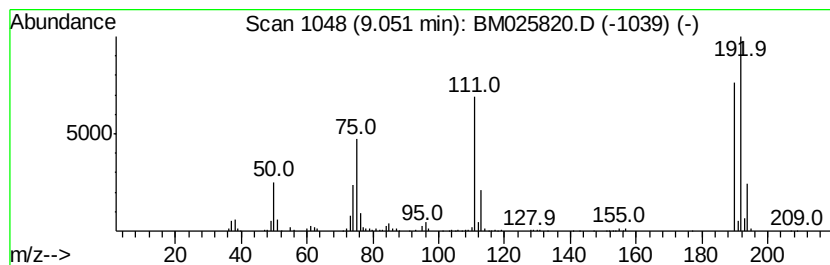
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 4-Chlorophenylselenol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	28.96 ng/ul	2045620	Napthalene-d8	10.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
2		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
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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Instrument :
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ClientSampleId :
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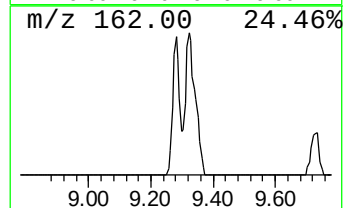
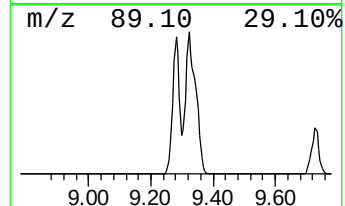
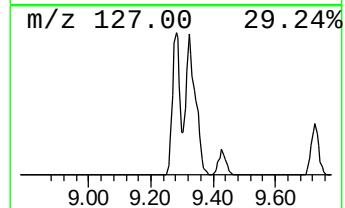
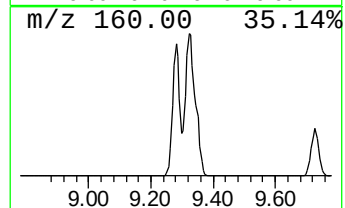
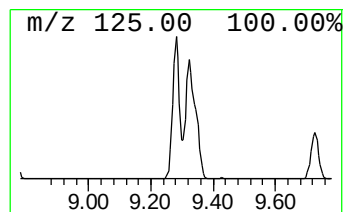
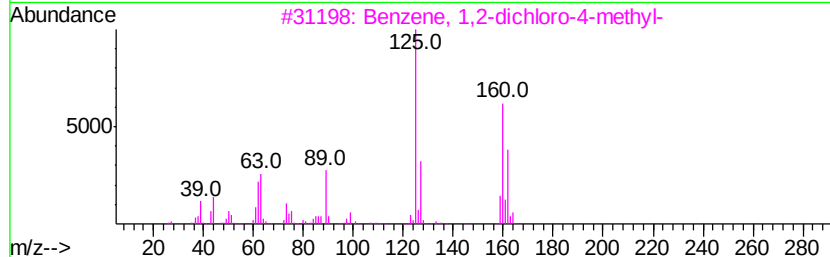
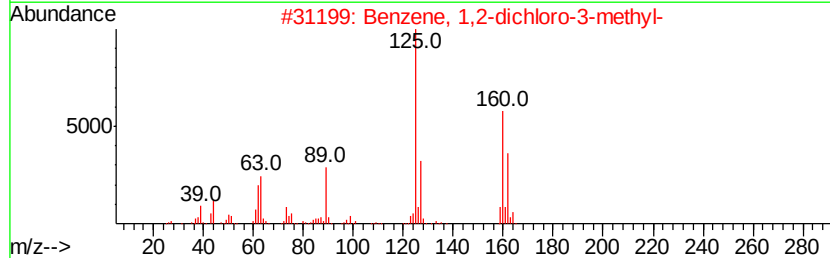
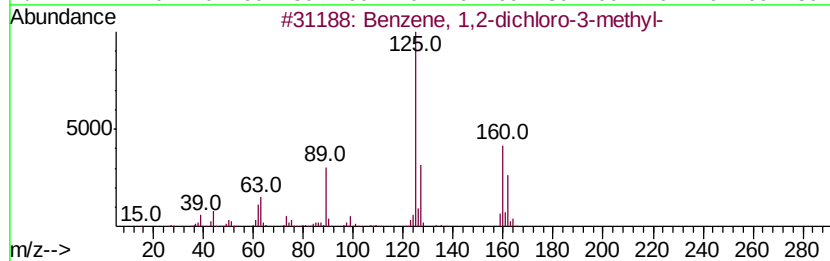
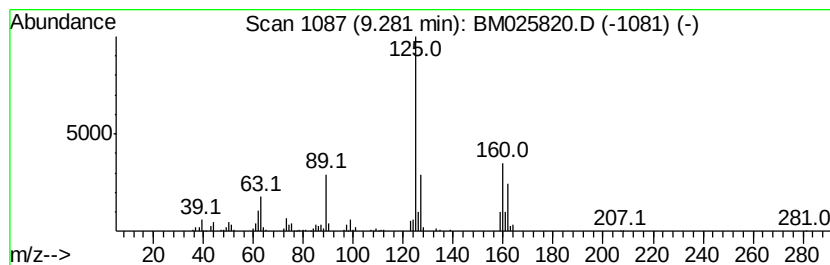
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Benzene, 1,2-dichloro-3-met... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.28	6.56 ng/ul	463503	Napthalene-d8	10.36

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



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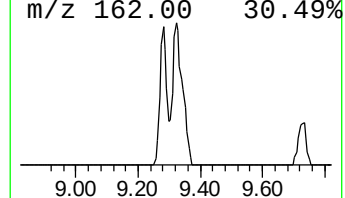
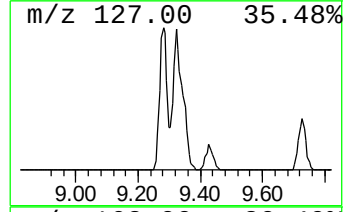
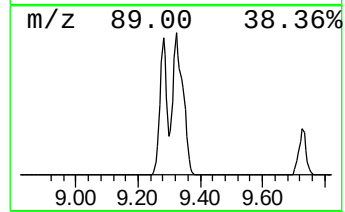
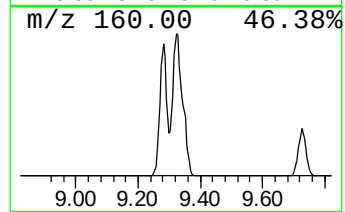
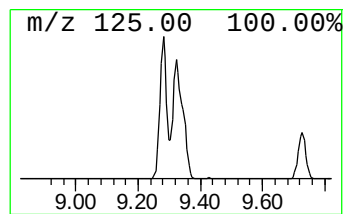
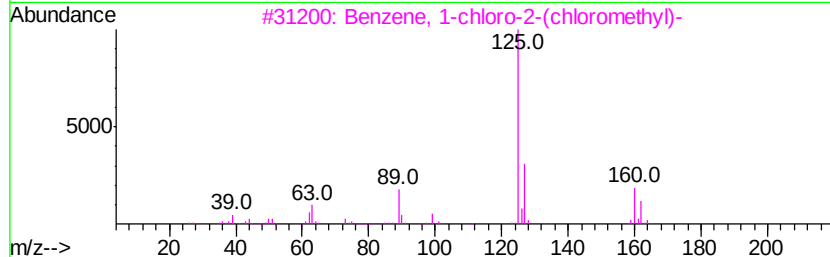
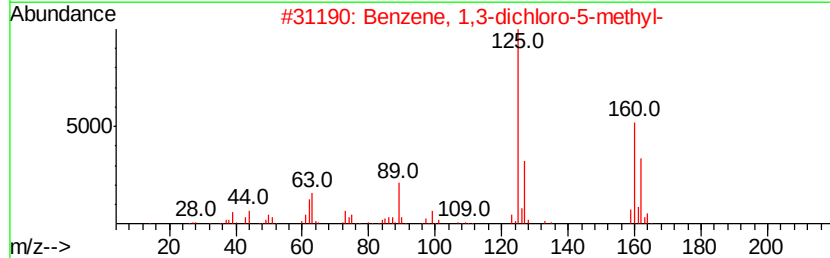
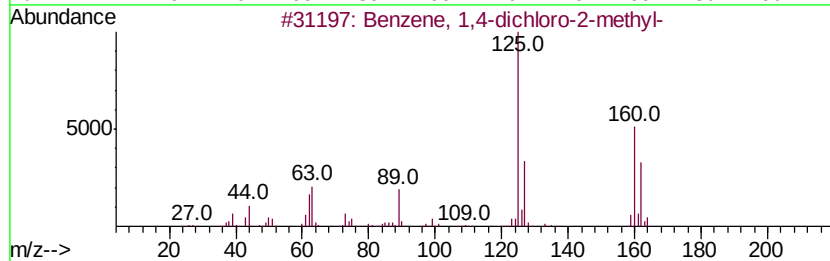
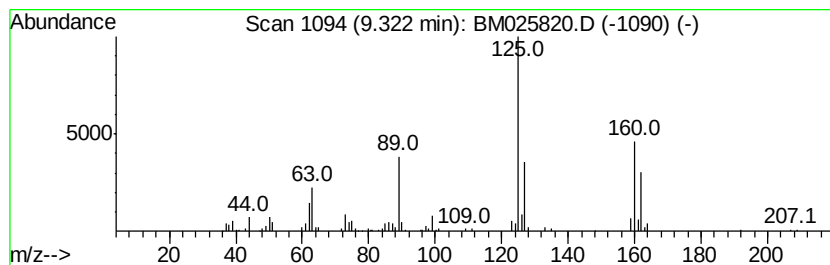
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,4-dichloro-2-met... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.32	6.16 ng/ul	435313	Napthalene-d8	10.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
2		Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	95
3		Benzene, 1-chloro-2-(chloromethyl)-	160	C7H6Cl2	000611-19-8	95
4		Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	94
5		Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	93



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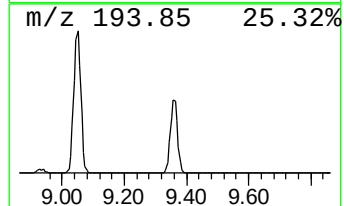
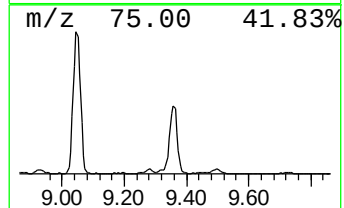
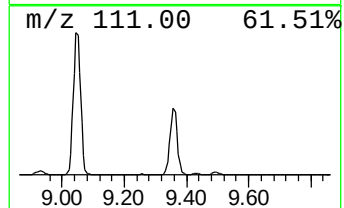
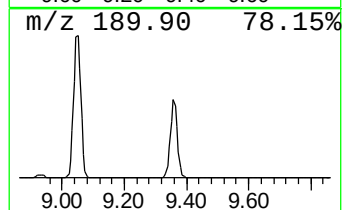
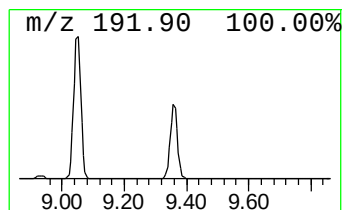
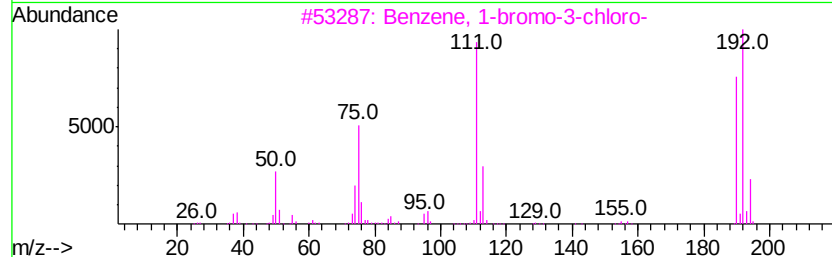
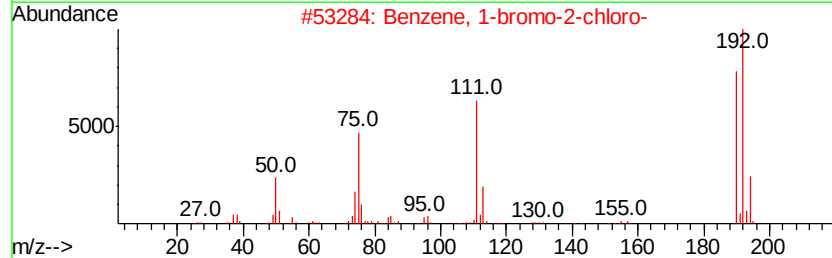
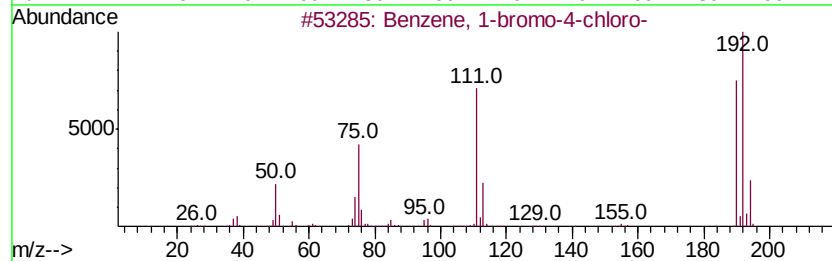
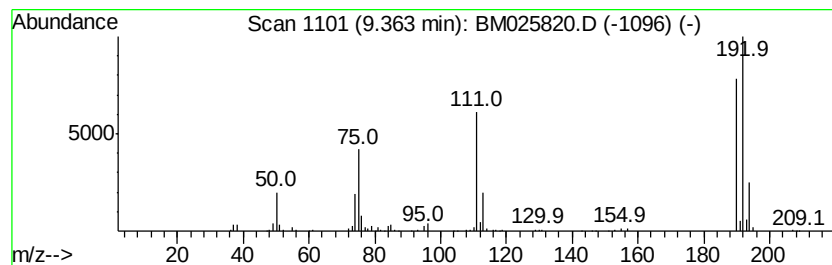
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzene, 1-bromo-4-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.36	17.07 ng/ul	1205610	Napthalene-d8	10.36

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	97
3		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
4		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	95
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	95



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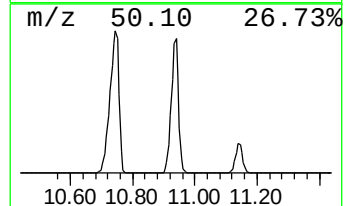
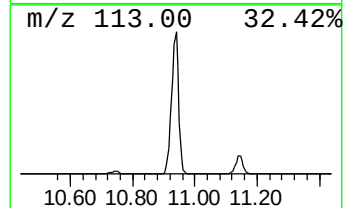
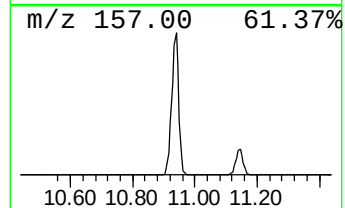
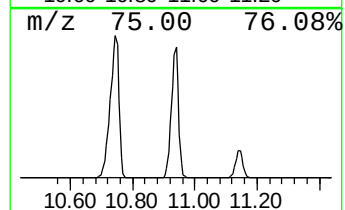
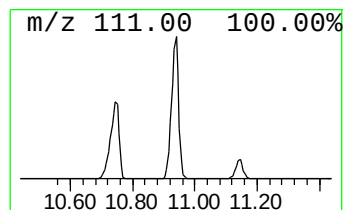
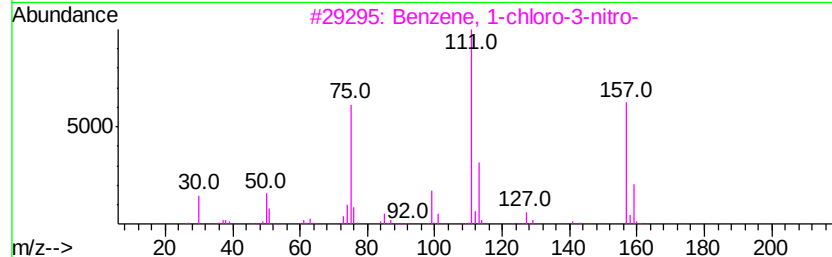
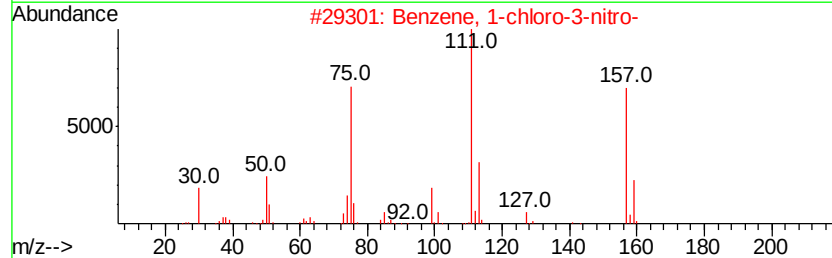
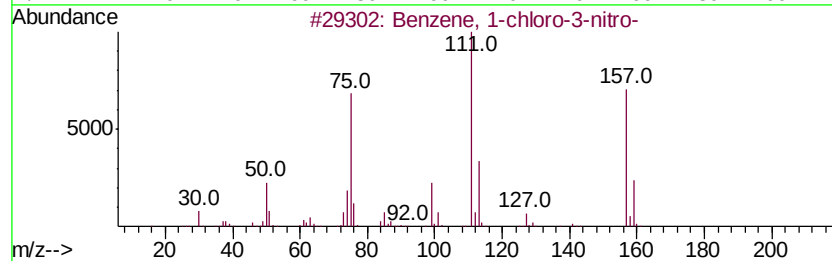
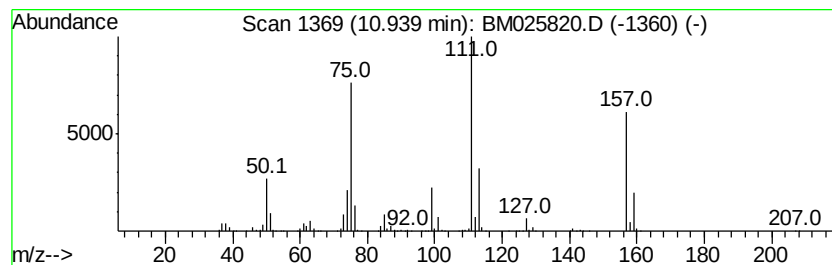
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	111.18 ng/ul	7852330	Naphthalene-d8	10.36

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	94
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025820.D
Acq On : 27 Mar 2020 03:56
Operator : CG/JU
Sample : L1968-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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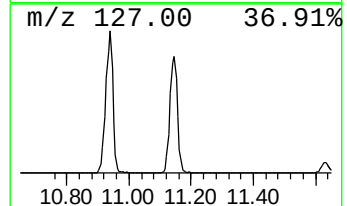
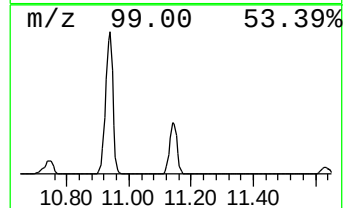
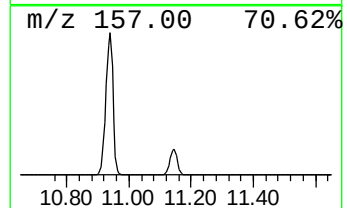
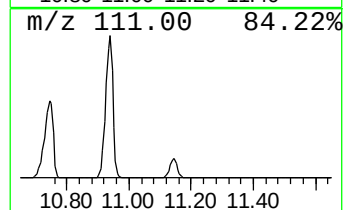
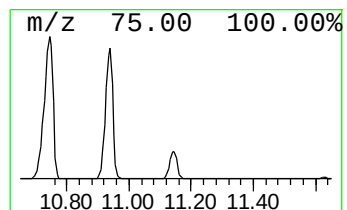
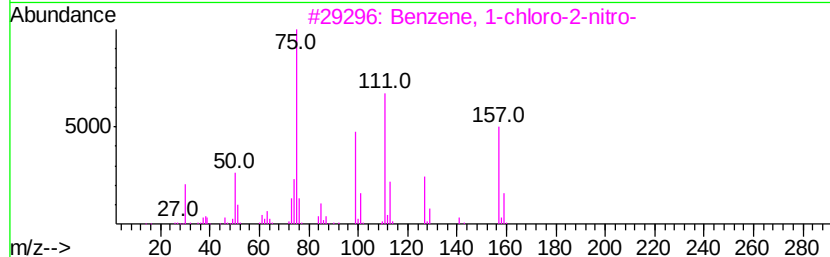
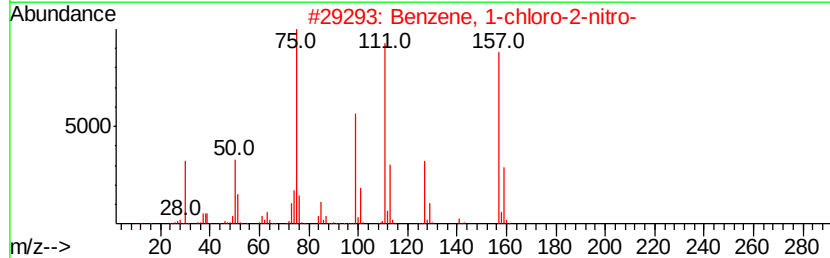
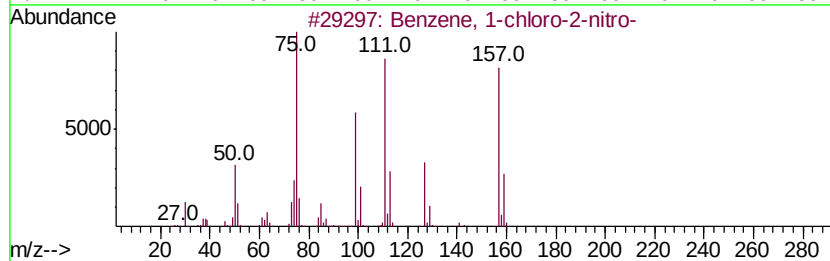
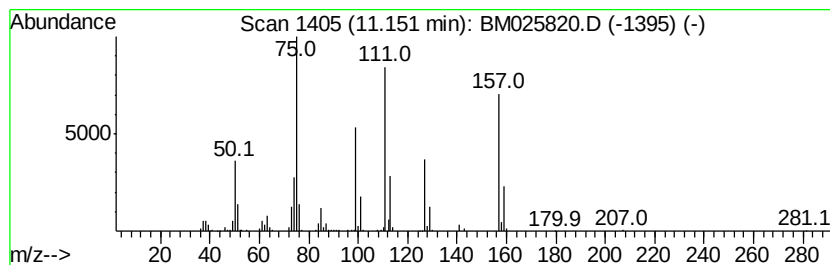
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzene, 1-chloro-2-nitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.15	23.65 ng/ul	1670170	Naphthalene-d8	10.36

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	97
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97



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Data File : BM025820.D
Acq On : 27 Mar 2020 03:56
Operator : CG/JU
Sample : L1968-16
Misc :
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ClientSampleID :
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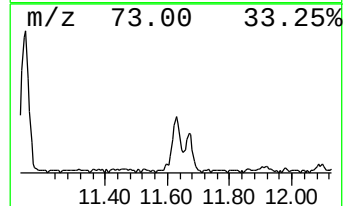
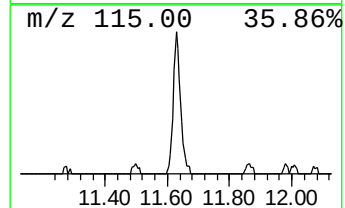
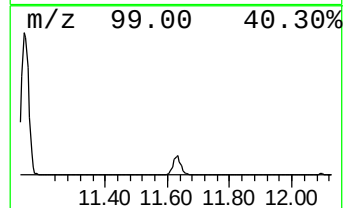
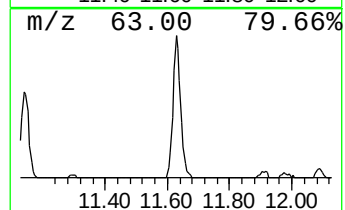
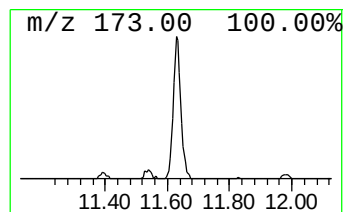
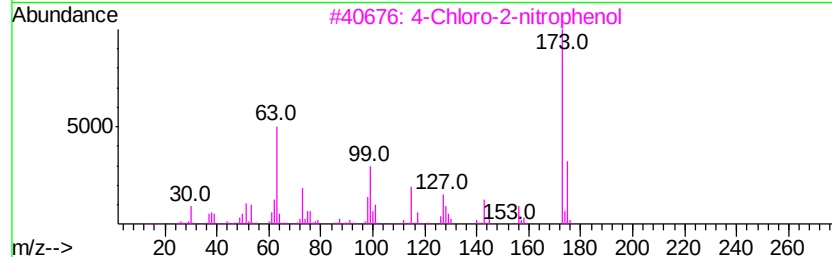
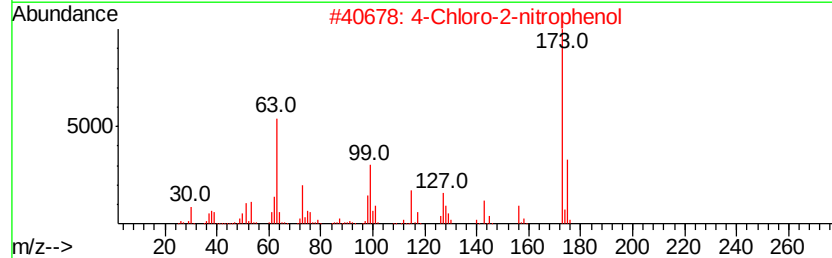
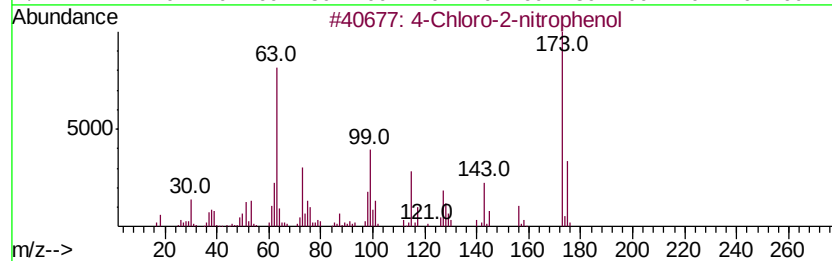
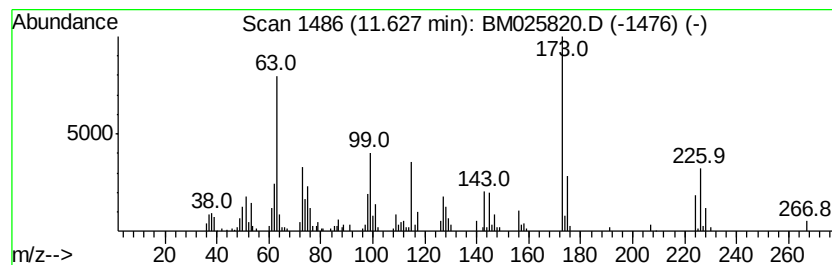
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.63	6.12 ng/ul	432317	Napthalene-d8	10.36

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	95
3			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	91
4			Phenol, 2-chloro-4-nitro-	173	C6H4ClNO3	000619-08-9	90
5			Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025820.D
Acq On : 27 Mar 2020 03:56
Operator : CG/JU
Sample : L1968-16
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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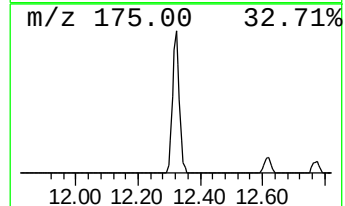
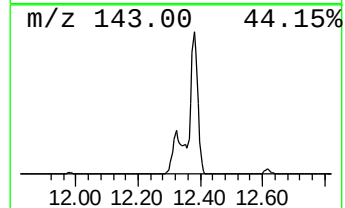
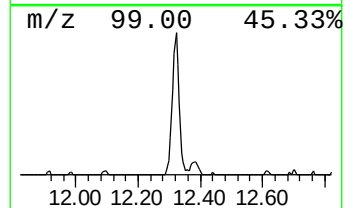
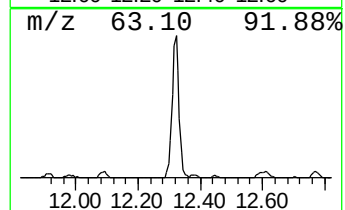
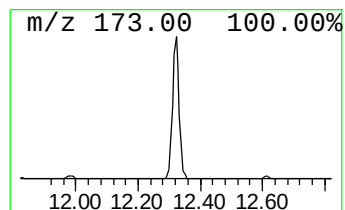
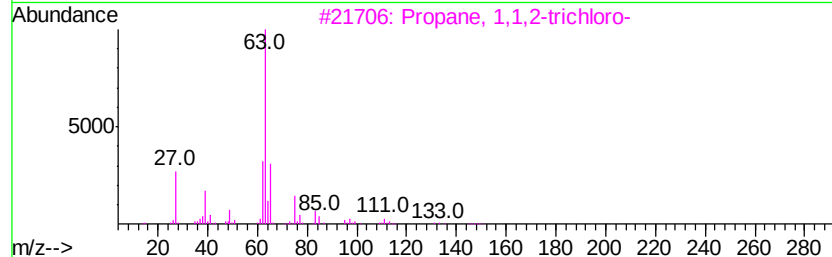
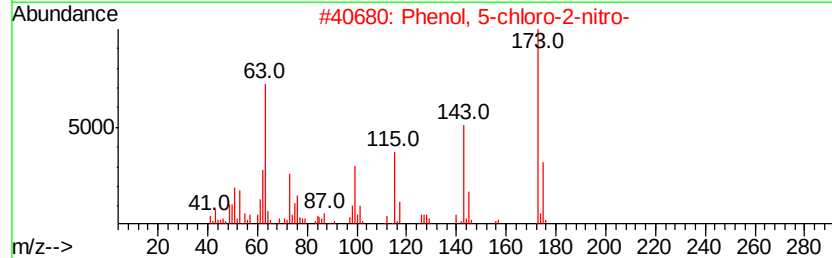
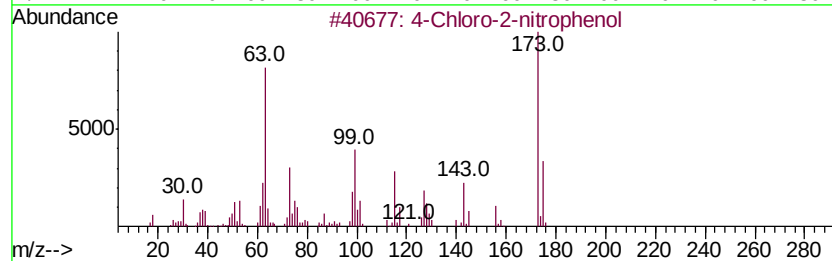
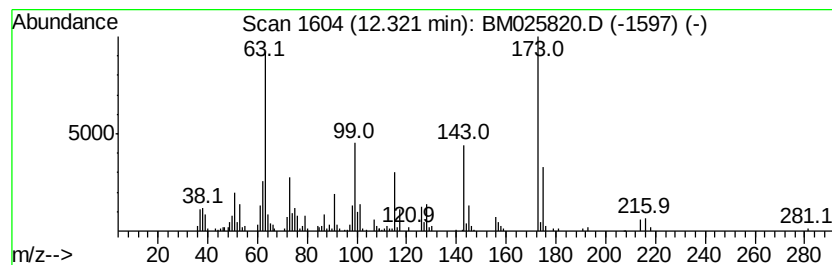
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Phenol, 5-chloro-2-nitro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	4.76 ng/ul	462480	Acenaphthene-d10	14.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	93
2			Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	83
3			Propane, 1,1,2-trichloro-	146	C3H5Cl3	000598-77-6	47
4			1-Bromo-1-chloroethane	142	C2H4BrCl	000593-96-4	43
5			Phenol, 2-chloro-4-nitro-	173	C6H4ClNO3	000619-08-9	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
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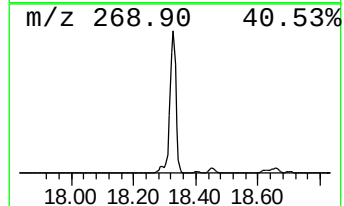
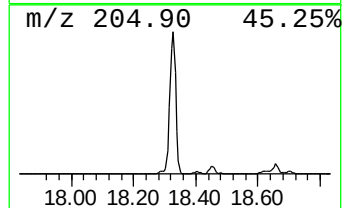
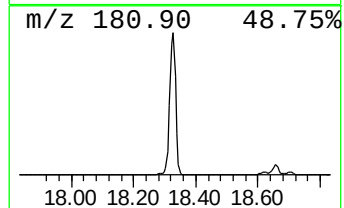
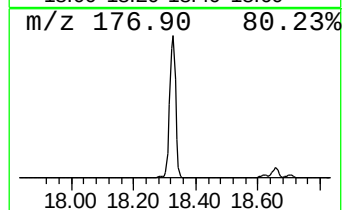
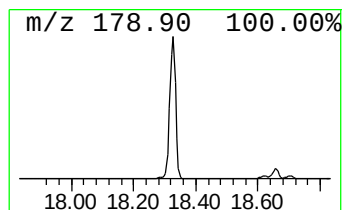
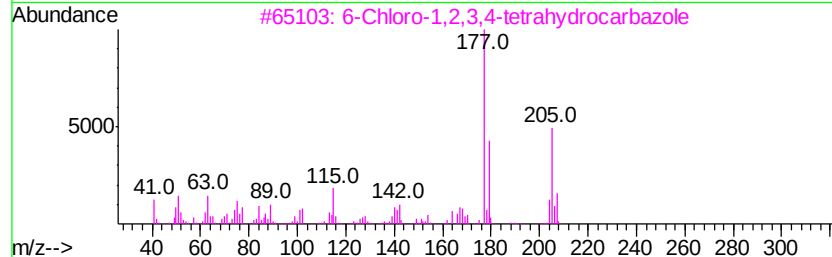
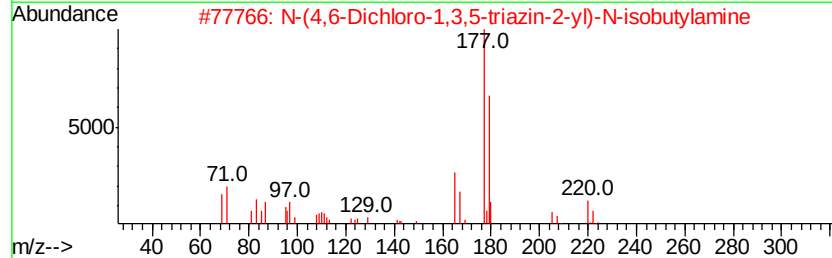
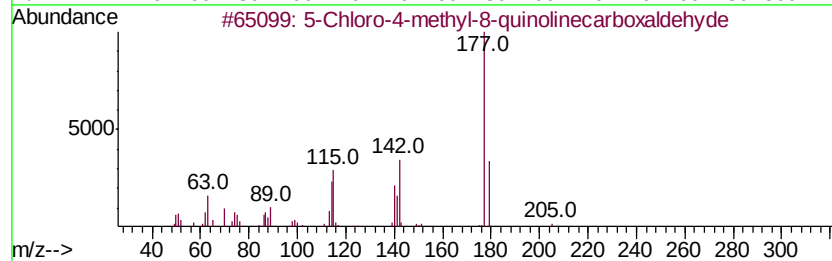
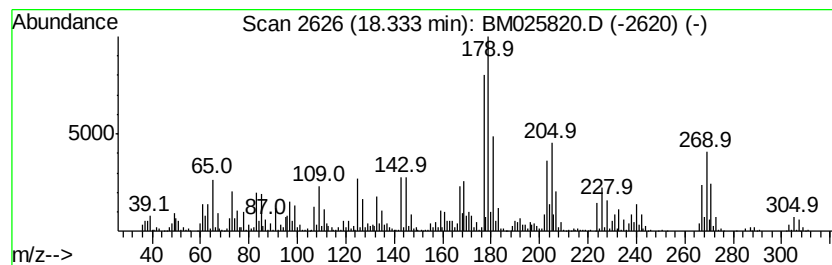
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.33	63.21 ng/ul	7140330	Phenanthrene-d10	16.95	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Chloro-4-methyl-8-quinolinecar...	205	C11H8ClNO	028662-47-7	27
2	N-(4,6-Dichloro-1,3,5-triazin-2-...	220	C7H10Cl2N4	1000326-88-2	27
3	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25
4	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	25
5	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025820.D
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Sample : L1968-16
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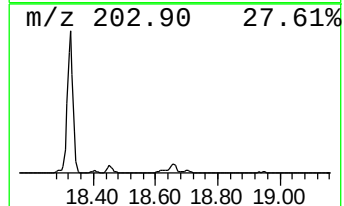
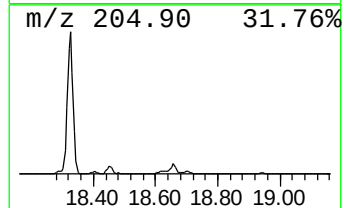
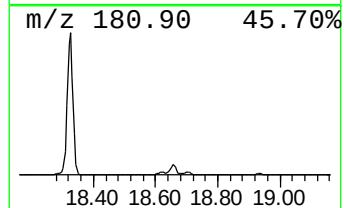
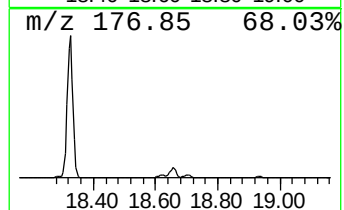
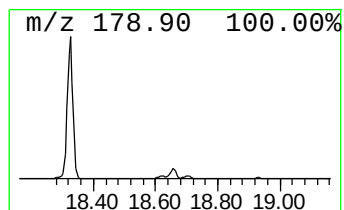
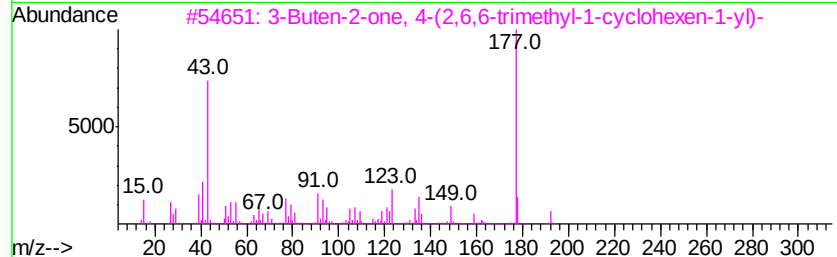
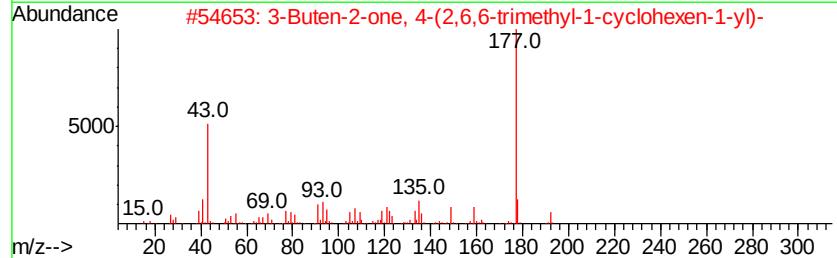
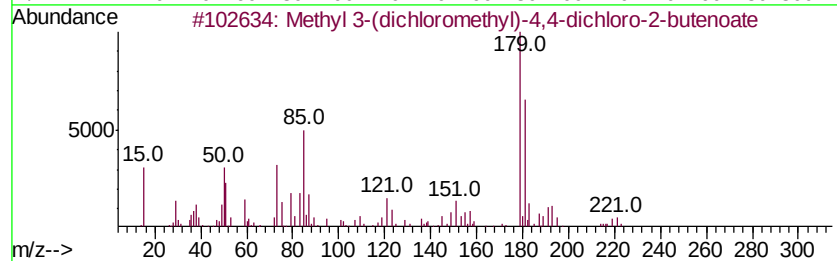
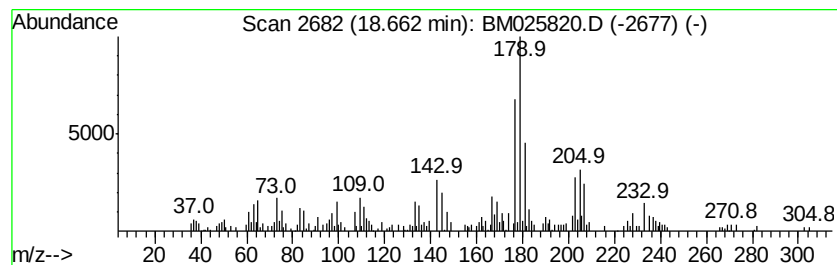
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-02 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.66	4.31 ng/ul	487285	Phenanthrene-d10	16.95
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Methyl 3-(dichloromethyl)-4,4-di...	250 C6H6Cl4O2	097055-33-9 12
2		3-Buten-2-one, 4-(2,6,6-trimethy...	192 C13H20O	014901-07-6 11
3		3-Buten-2-one, 4-(2,6,6-trimethy...	192 C13H20O	014901-07-6 10
4		3-Buten-2-one, 4-(2,6,6-trimethy...	192 C13H20O	014901-07-6 10
5		3,4-Dihydroxy-1,6-bis-(3-methoxy...	354 C20H18O6	1000297-24-6 10



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025820.D
Acq On : 27 Mar 2020 03:56
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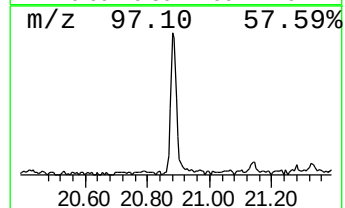
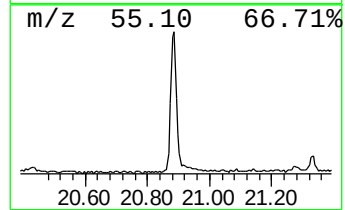
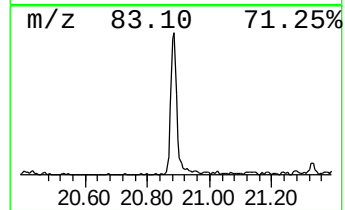
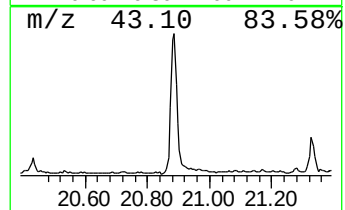
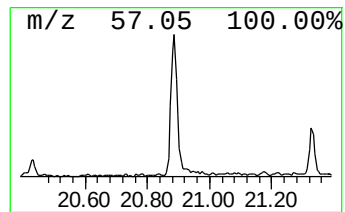
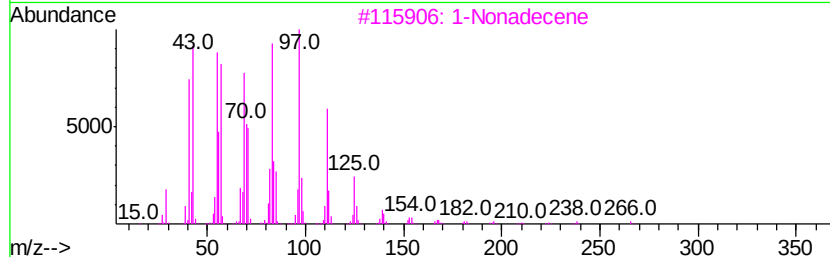
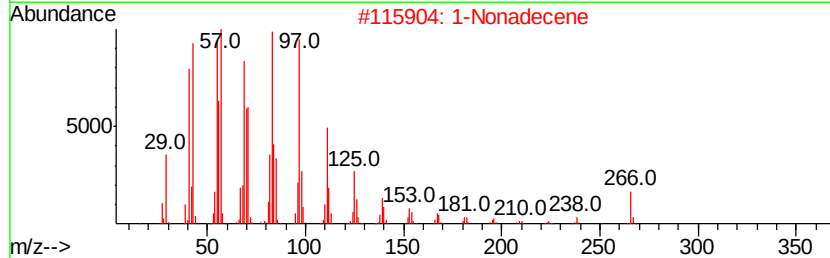
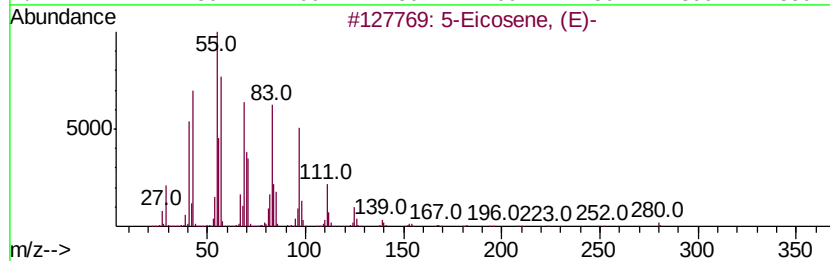
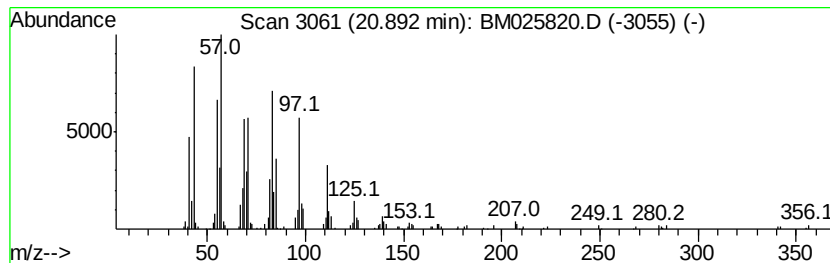
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.89	3.61 ng/ul	493784	Chrysene-d12	21.14

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Eicosene, (E)-	280	C20H40	074685-30-6	98
2			1-Nonadecene	266	C19H38	018435-45-5	96
3			1-Nonadecene	266	C19H38	018435-45-5	95
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5			n-Tetracosanol-1	354	C24H50O	000506-51-4	94



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032620\
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Acq On : 27 Mar 2020 03:56
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Misc :
ALS Vial : 19 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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
4-Chlorophenylsel...	9.05	29.0	ng/ul	2045620	2	10.36	1412530	20.0
Benzene, 1,2-dich...	9.28	6.6	ng/ul	463503	2	10.36	1412530	20.0
Benzene, 1,4-dich...	9.32	6.2	ng/ul	435313	2	10.36	1412530	20.0
Benzene, 1-bromo-...	9.36	17.1	ng/ul	1205610	2	10.36	1412530	20.0
Benzene, 1-chloro...	10.94	111.2	ng/ul	7852330	2	10.36	1412530	20.0
Benzene, 1-chloro...	11.15	23.6	ng/ul	1670170	2	10.36	1412530	20.0
4-Chloro-2-nitrop...	11.63	6.1	ng/ul	432317	2	10.36	1412530	20.0
Phenol, 5-chloro-...	12.32	4.8	ng/ul	462480	3	14.20	1941860	20.0
unknown-01	18.33	63.2	ng/ul	7140330	4	16.95	2259100	20.0
unknown-02	18.66	4.3	ng/ul	487285	4	16.95	2259100	20.0
(DEL) Alkane: Str...	20.89	3.6	ng/ul	493784	5	21.14	2736740	20.0