

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
 Data File : BM025838.D
 Acq On : 27 Mar 2020 15:15
 Operator : CG/JU
 Sample : L2005-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0DF0

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	4.993	340	358	372	RBV7	67784117	570517272	100.00%	31.047%
2	6.269	568	575	586	RBV	382090	622413	0.11%	0.034%
3	6.951	679	691	700	RBV2	336780	1087008	0.19%	0.059%
4	7.116	712	719	737	rVB2	529773	1720018	0.30%	0.094%
5	7.481	770	781	797	RBV	13619515	67814113	11.89%	3.690%
6	7.681	797	815	825	RBV8	72216936	523992967	91.85%	28.515%
7	8.040	853	876	883	RBV8	72338656	526022026	92.20%	28.625%
8	8.287	912	918	926	rVB	609776	934706	0.16%	0.051%
9	8.722	985	992	996	RBV	900378	1428374	0.25%	0.078%
10	9.051	1041	1048	1059	rVB	500695	783938	0.14%	0.043%
11	9.381	1095	1104	1109	RBV	3558231	5902134	1.03%	0.321%
12	9.434	1109	1113	1121	rVB	771063	1164829	0.20%	0.063%
13	9.963	1197	1203	1210	RBV	1093323	1804550	0.32%	0.098%
14	10.034	1210	1215	1223	rVB	395364	656244	0.12%	0.036%
15	10.128	1223	1231	1237	RBV	668160	1222186	0.21%	0.067%
16	10.228	1237	1248	1258	RBV	34433761	62838436	11.01%	3.420%
17	10.345	1261	1268	1275	rVB	985736	1550812	0.27%	0.084%
18	10.457	1278	1287	1293	RBV	5322600	9091708	1.59%	0.495%
19	10.728	1323	1333	1340	RBV	8824357	14202447	2.49%	0.773%
20	10.928	1360	1367	1376	RBV	584252	927450	0.16%	0.050%
21	12.004	1542	1550	1562	RBV	349403	643223	0.11%	0.035%
22	12.386	1608	1615	1624	rVB2	848320	1354090	0.24%	0.074%
23	12.622	1637	1655	1664	rVB2	349374	637115	0.11%	0.035%
24	13.622	1818	1825	1830	rVV	1963639	2602285	0.46%	0.142%
25	13.892	1865	1871	1880	rVB	2374857	3333077	0.58%	0.181%
26	14.204	1918	1924	1930	rVB	1424048	2044100	0.36%	0.111%
27	15.204	2087	2094	2107	rVB	2824244	4046216	0.71%	0.220%
28	15.321	2109	2114	2124	RBV	828927	1193425	0.21%	0.065%
29	16.951	2385	2391	2400	rVB	1759230	2438029	0.43%	0.133%
30	17.045	2401	2407	2420	RBV	3228031	4455340	0.78%	0.242%
31	19.345	2792	2798	2804	RBV	4175561	5336394	0.94%	0.290%
32	21.139	3098	3103	3111	RBV	2498096	3078798	0.54%	0.168%
33	23.156	3439	3446	3453	RBV	3715569	6194017	1.09%	0.337%
34	23.286	3462	3468	3478	rVB2	1824672	3360469	0.59%	0.183%

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35	24.080	3597	3603	3611	rVB	362216	700306	0.12%	0.038%
36	25.580	3851	3858	3867	rVB	531932	1273239	0.22%	0.069%
37	27.603	4195	4202	4213	rVB3	208099	642745	0.11%	0.035%

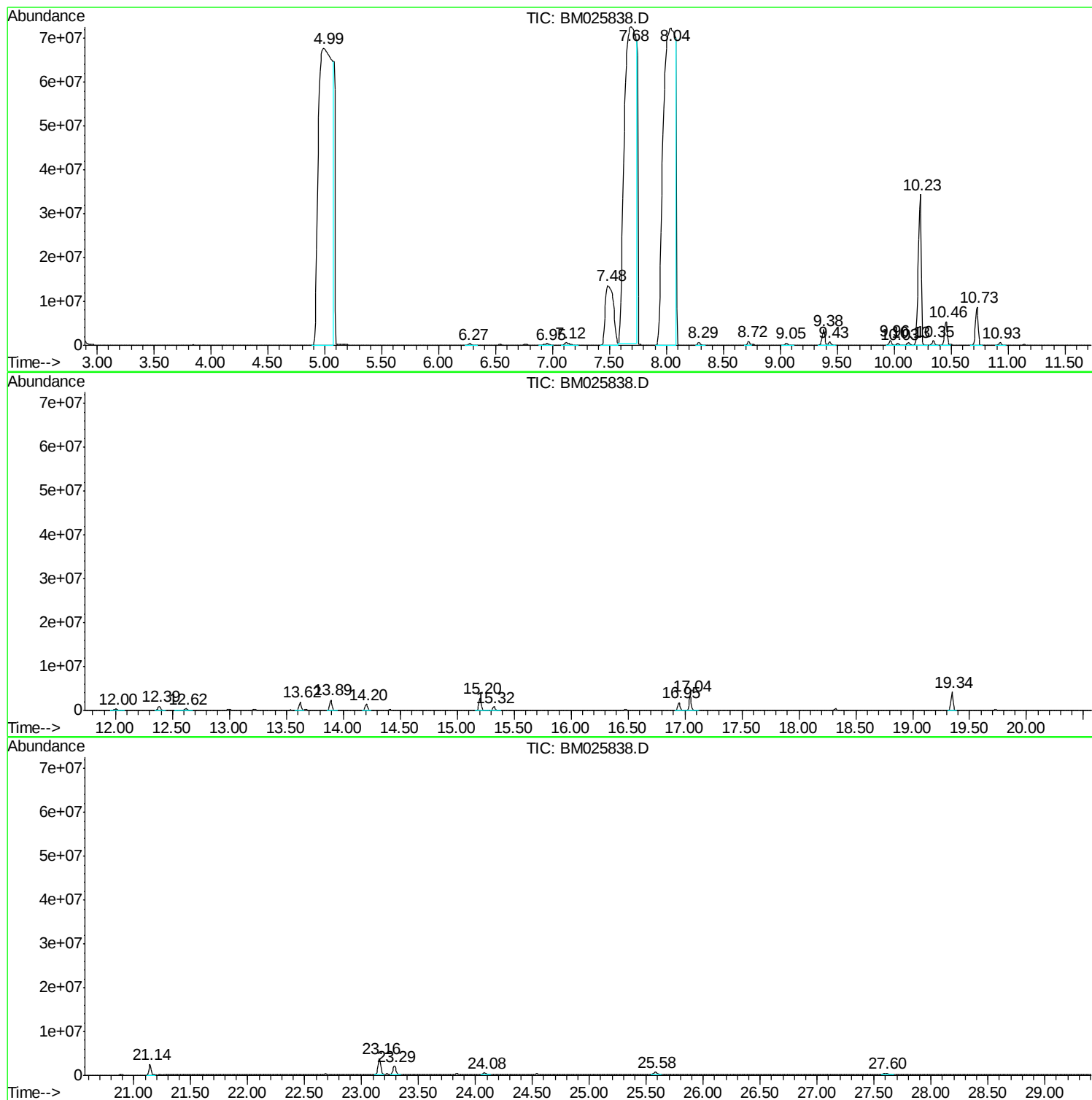
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Quant Title : SVOA CALIBRATION

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



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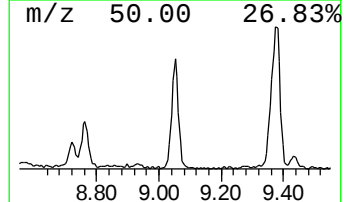
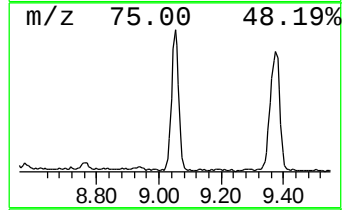
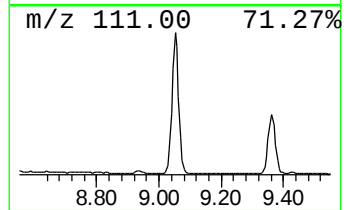
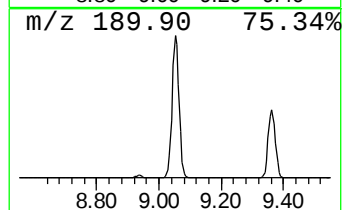
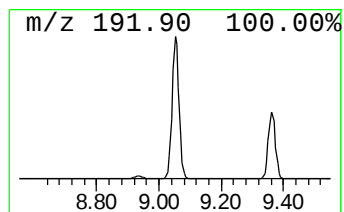
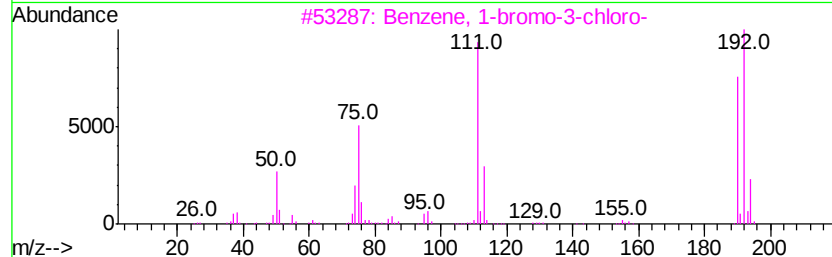
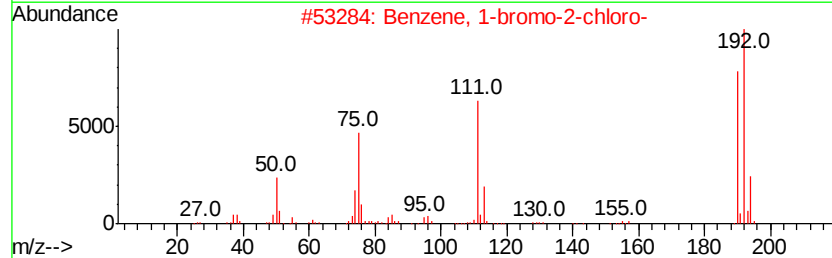
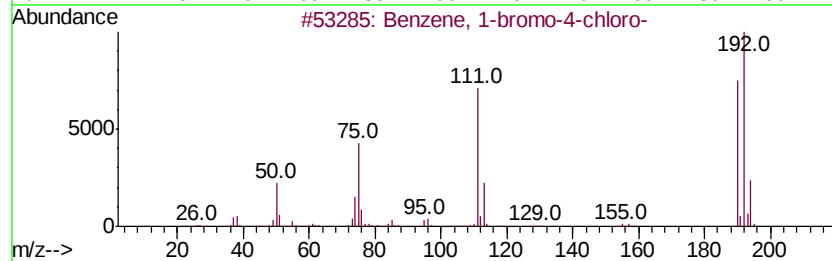
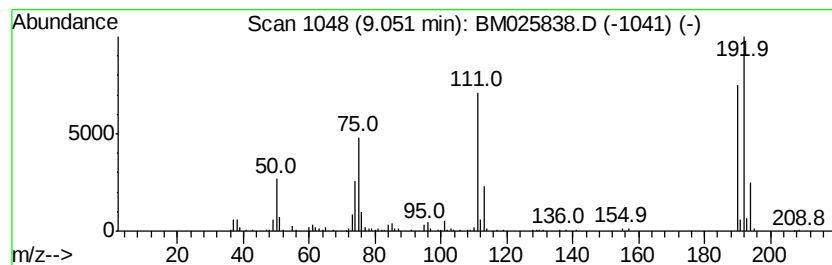
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	10.11 ng/ul	783938	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		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	97



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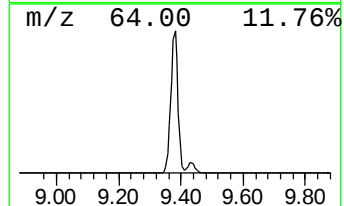
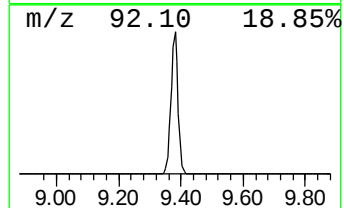
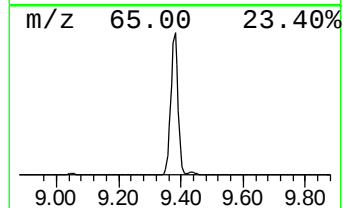
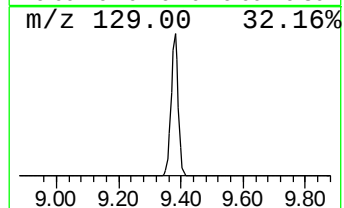
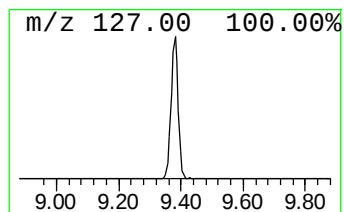
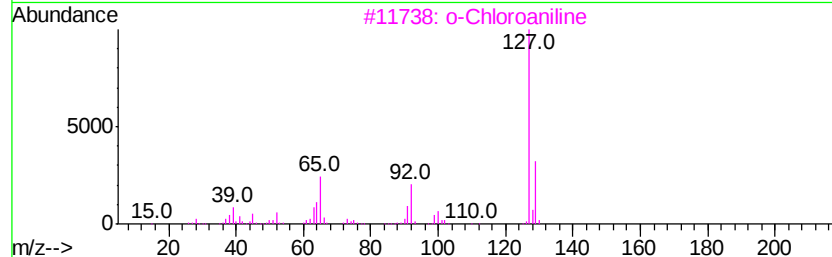
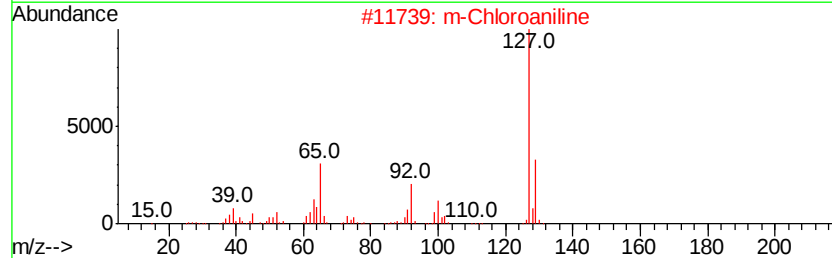
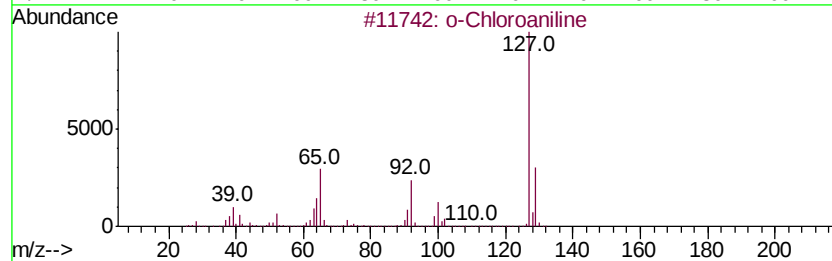
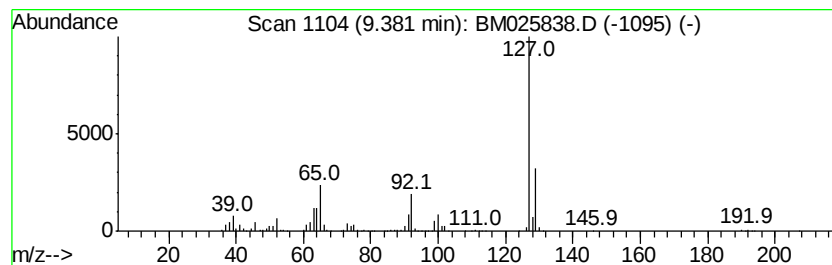
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM031420MA.M
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 5 o-Chloroaniline Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.38	76.12 ng/ul	5902130	Naphthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Chloroaniline	127	C6H6ClN	000095-51-2	97
2		m-Chloroaniline	127	C6H6ClN	000108-42-9	96
3		o-Chloroaniline	127	C6H6ClN	000095-51-2	96
4		m-Chloroaniline	127	C6H6ClN	000108-42-9	96
5		m-Chloroaniline	127	C6H6ClN	000108-42-9	94



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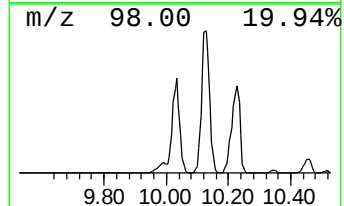
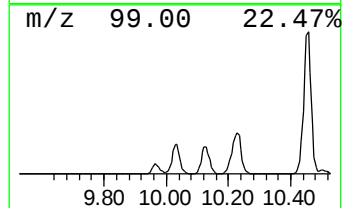
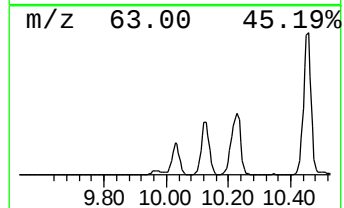
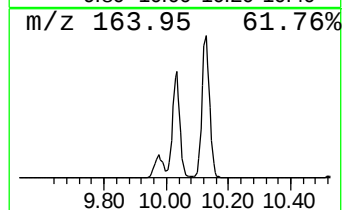
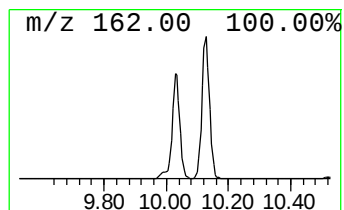
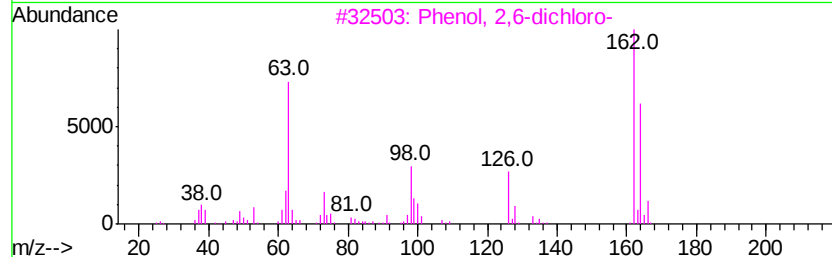
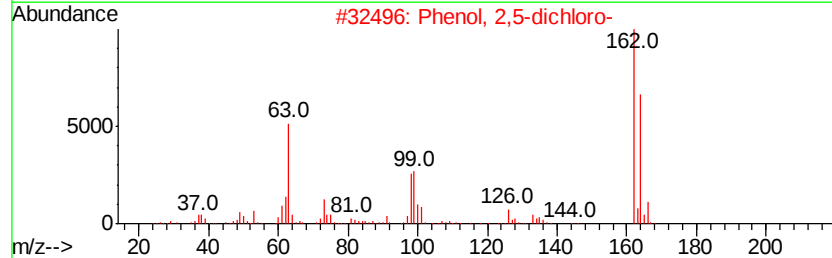
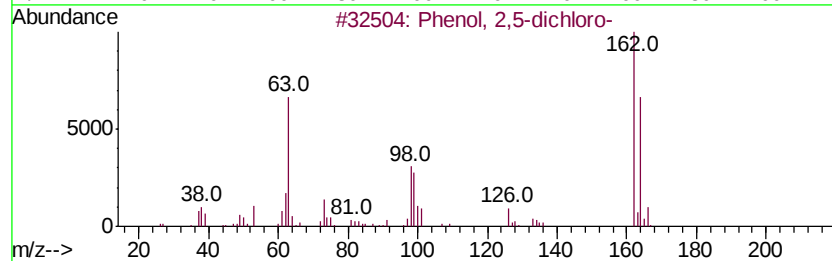
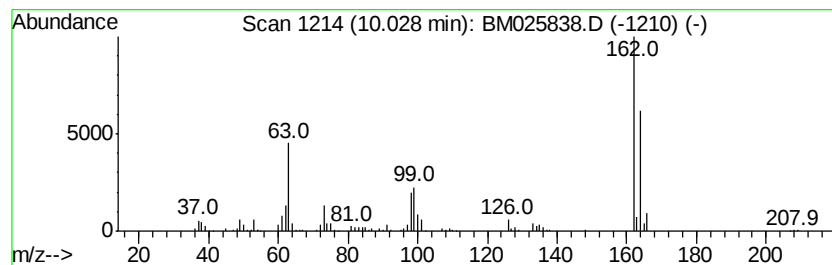
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 6 Phenol, 2,5-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.03	8.46 ng/ul	656244	Napthalene-d8	10.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Phenol, 2,5-dichloro-	162 C6H4Cl2O	000583-78-8	97
2	Phenol, 2,5-dichloro-	162 C6H4Cl2O	000583-78-8	94
3	Phenol, 2,6-dichloro-	162 C6H4Cl2O	000087-65-0	94
4	Phenol, 2,4-dichloro-	162 C6H4Cl2O	000120-83-2	94
5	Phenol, 2,6-dichloro-	162 C6H4Cl2O	000087-65-0	94



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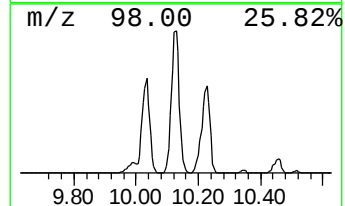
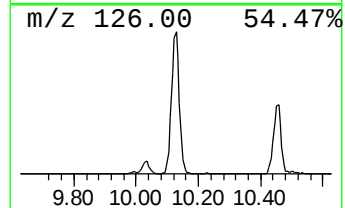
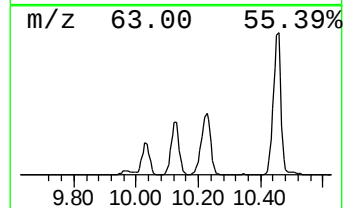
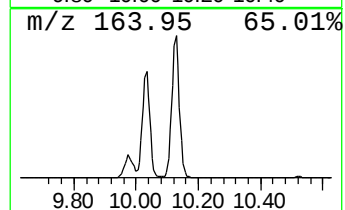
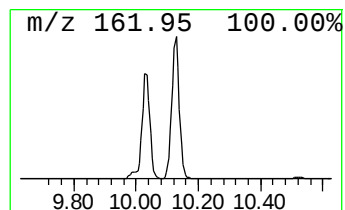
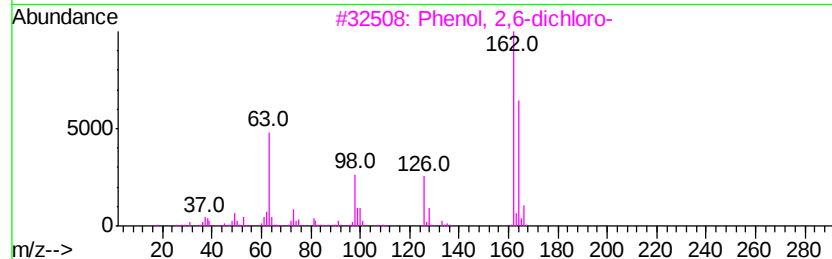
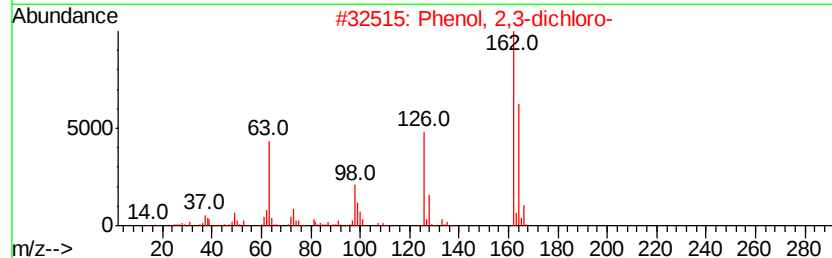
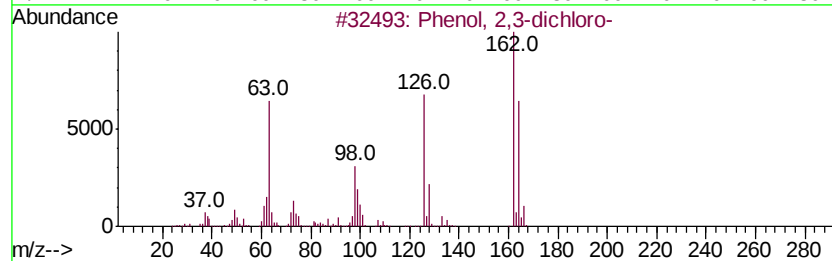
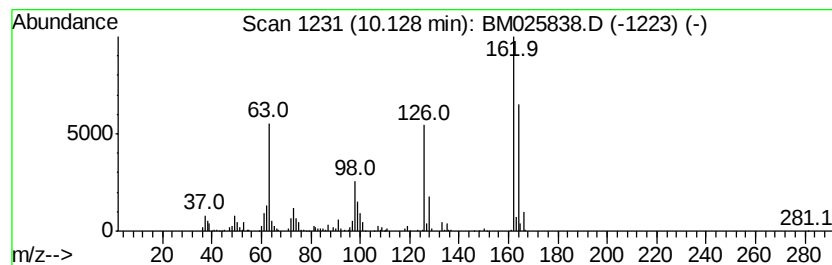
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Peak Number 7 Phenol, 2,3-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.13	15.76 ng/ul	1222190	Napthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	98
2		Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	98
3		Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	97
4		Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	96
5		Phenol, 2,4-dichloro-	162	C6H4Cl2O	000120-83-2	95



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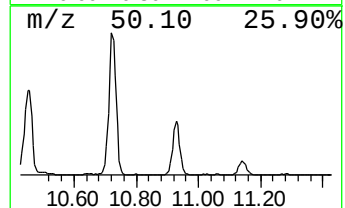
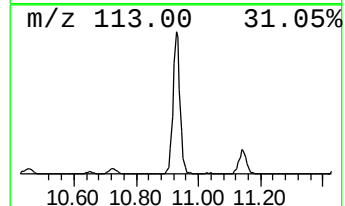
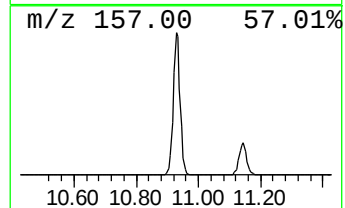
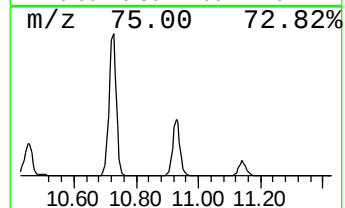
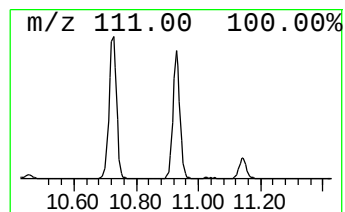
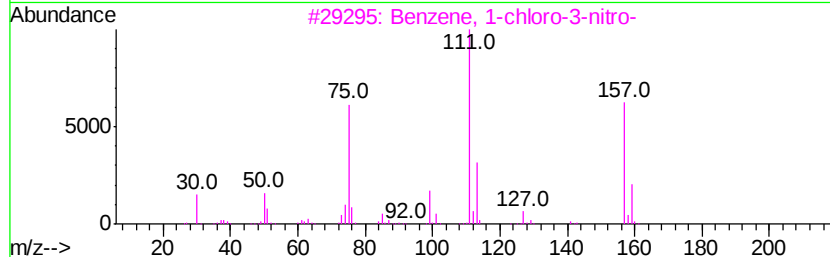
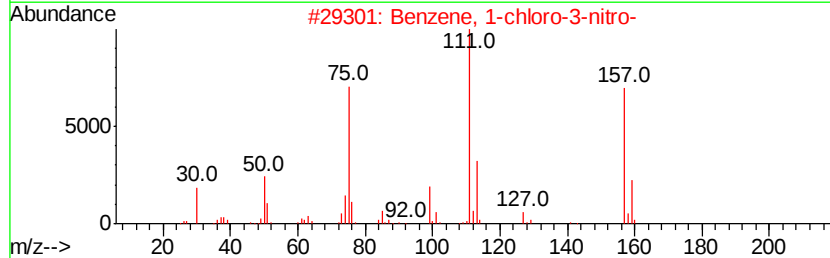
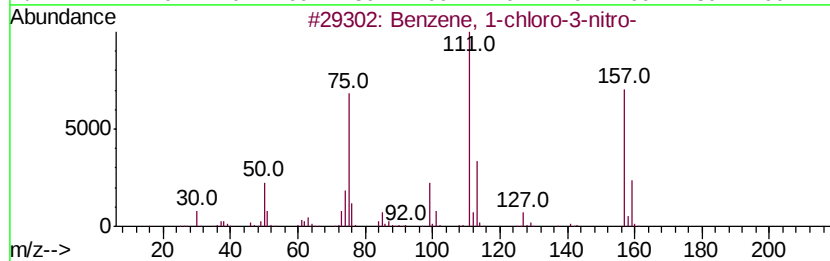
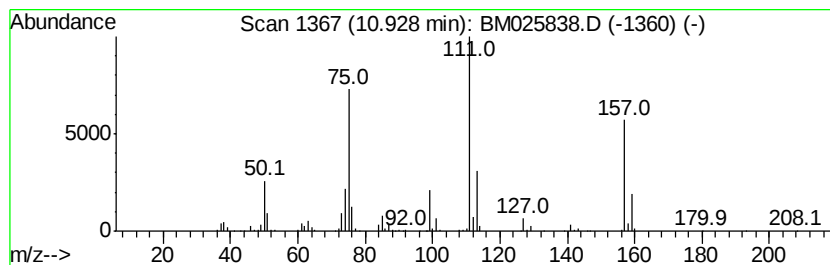
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TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-chloro-3-nitro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.93	11.96 ng/ul	927450	Naphthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025838.D
Acq On : 27 Mar 2020 15:15
Operator : CG/JU
Sample : L2005-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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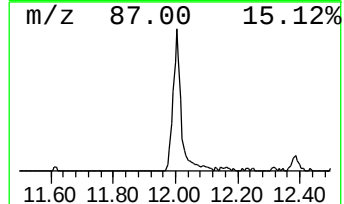
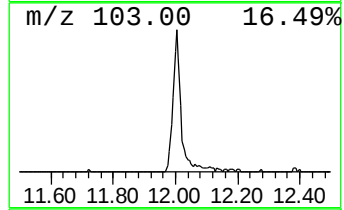
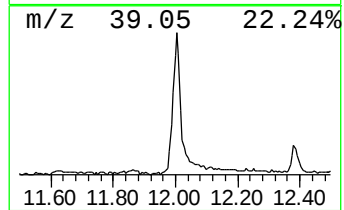
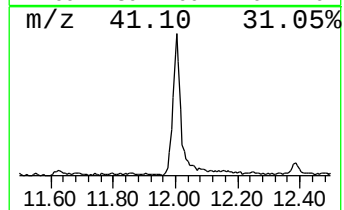
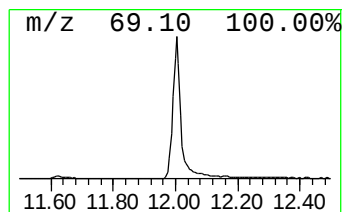
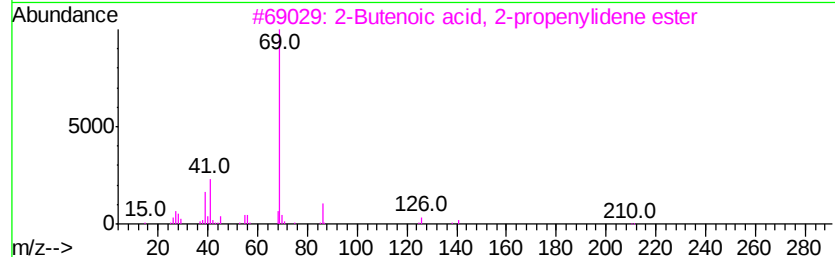
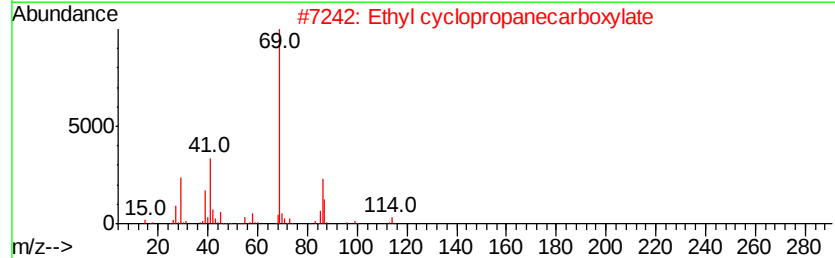
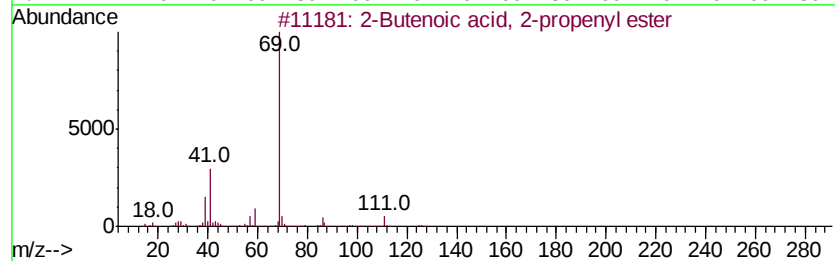
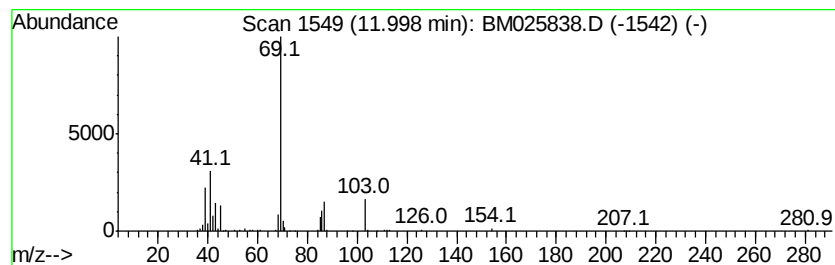
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 2-Butenoic acid, 2-propenyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	8.30 ng/ul	643223	Naphthalene-d8	10.35

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Butenoic acid, 2-propenyl ester	126	C7H10O2	020474-93-5	53
2		Ethyl cyclopropanecarboxylate	114	C6H10O2	004606-07-9	45
3		2-Butenoic acid, 2-propenylidene...	210	C11H14O4	055030-70-1	42
4		Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	33
5		2-Propenoic acid, 2-methyl-, 2-p...	126	C7H10O2	000096-05-9	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025838.D
Acq On : 27 Mar 2020 15:15
Operator : CG/JU
Sample : L2005-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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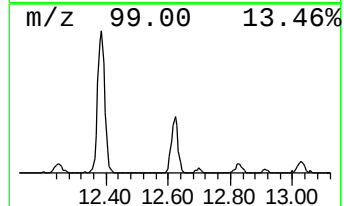
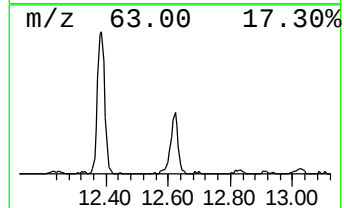
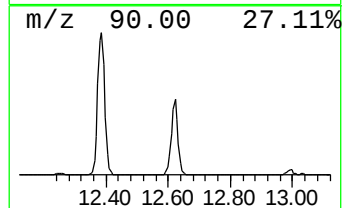
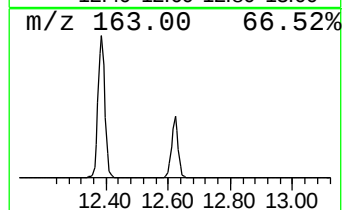
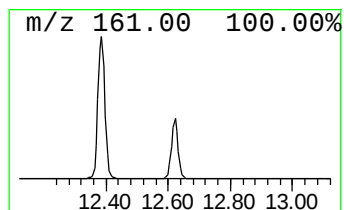
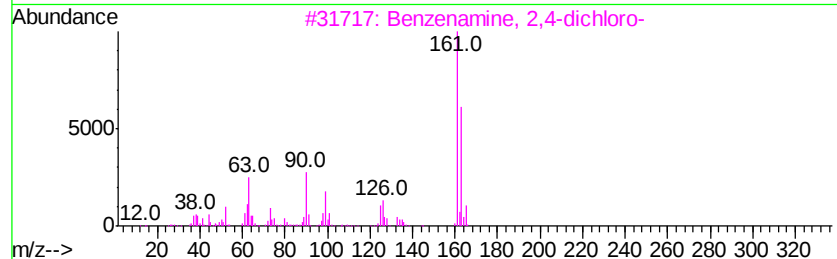
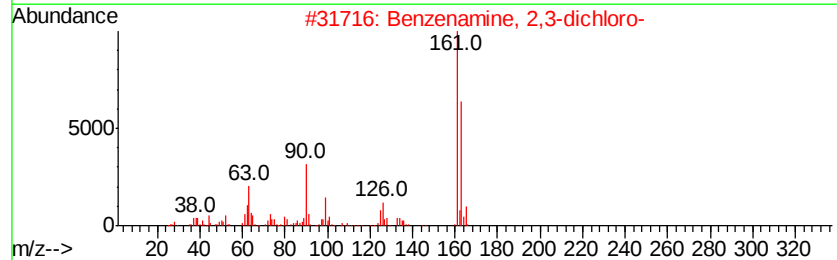
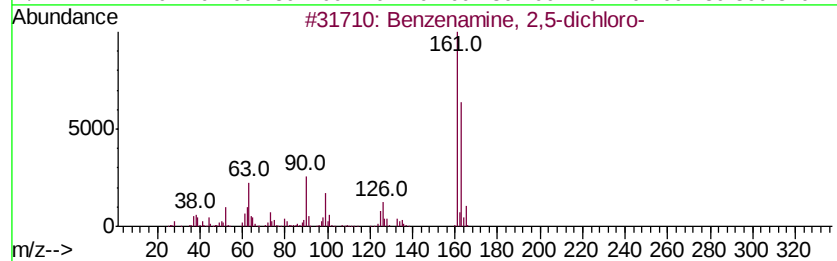
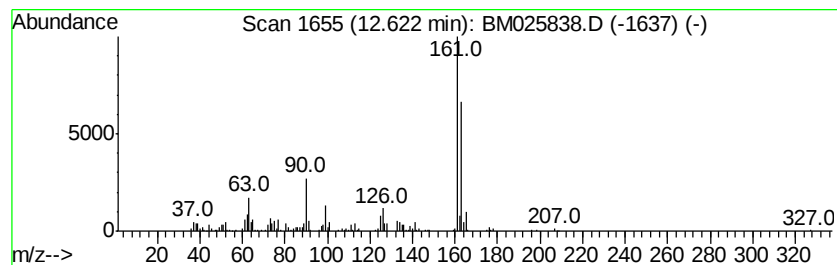
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Benzenamine, 2,5-dichloro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.62	6.23 ng/ul	637115	Acenaphthene-d10	14.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	98
2			Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	98
3			Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	96
4			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	96
5			Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025838.D
Acq On : 27 Mar 2020 15:15
Operator : CG/JU
Sample : L2005-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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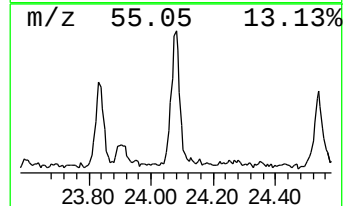
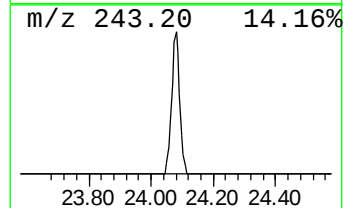
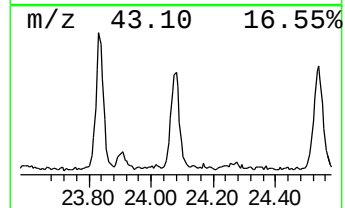
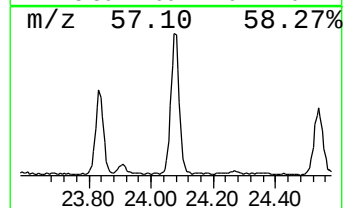
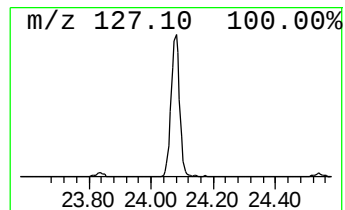
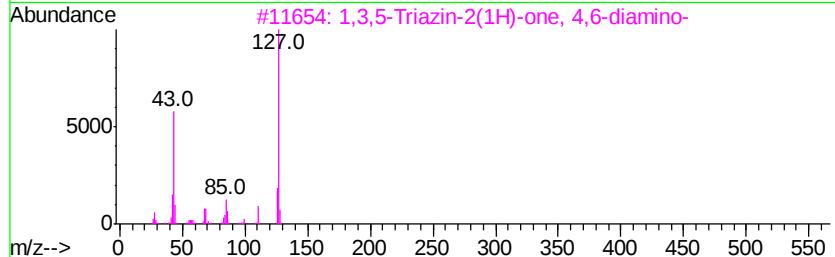
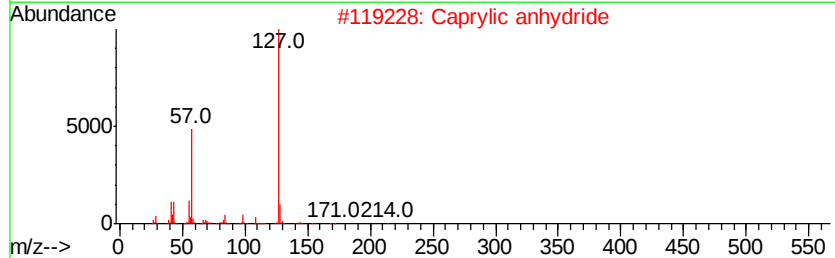
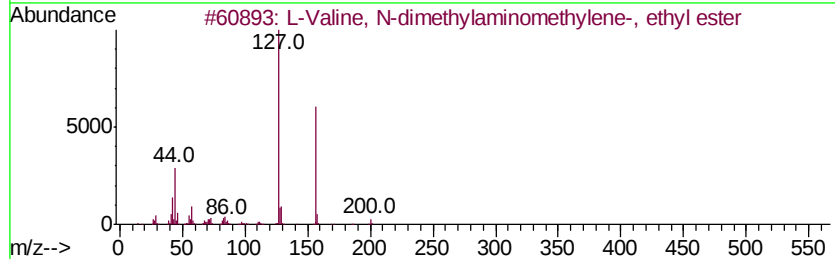
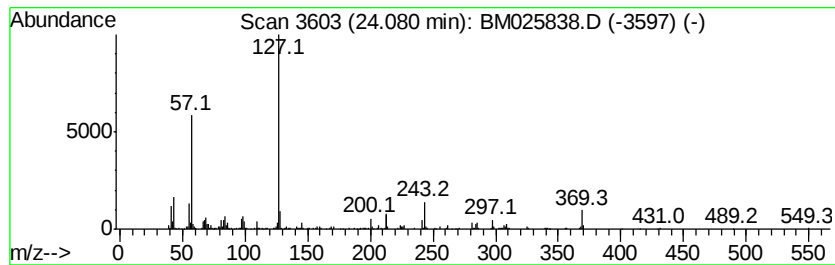
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 L-Valine, N-dimethylaminome... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	4.17 ng/ul	700306	Perylene-d12	23.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	L-Valine, N-dimethylaminomethyle...	200	C10H20N2O2	1000375-63-8	53
2		Caprylic anhydride	270	C16H30O3	000623-66-5	53
3		1,3,5-Triazin-2(1H)-one, 4,6-dia...	127	C3H5N5O	000645-92-1	53
4		Hexanoic acid, 2-ethyl-, anhydride	270	C16H30O3	036765-89-6	53
5		(Chlorophenoxymethylenimino)sulf...	281	C7H5ClF5NOS	081439-21-6	47



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025838.D
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Operator : CG/JU
Sample : L2005-08
Misc :
ALS Vial : 35 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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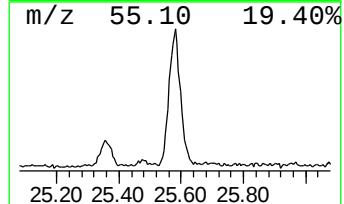
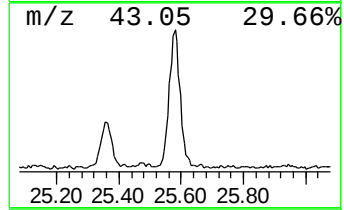
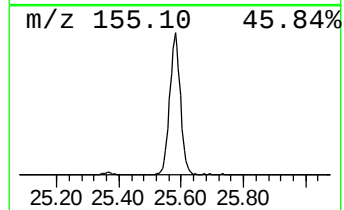
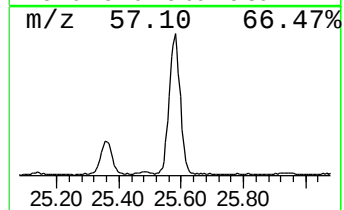
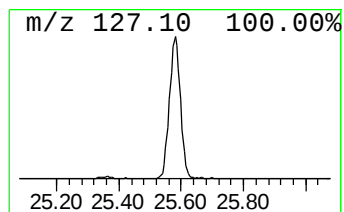
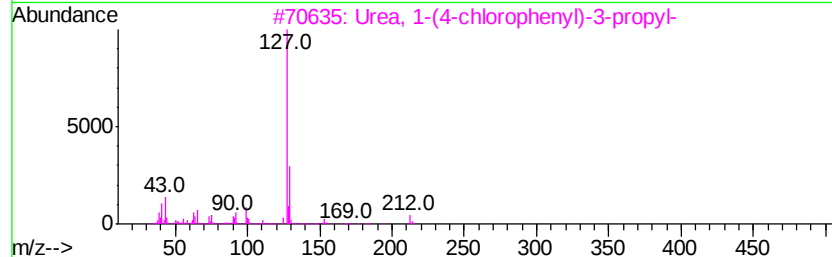
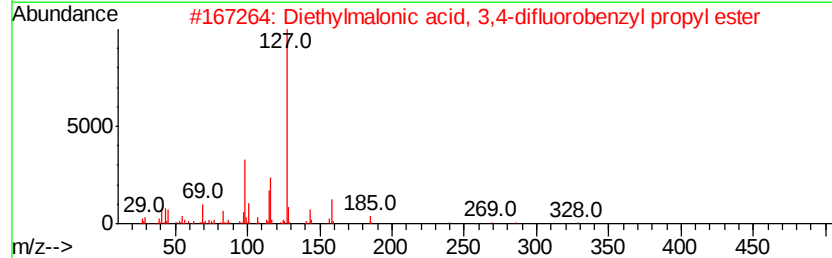
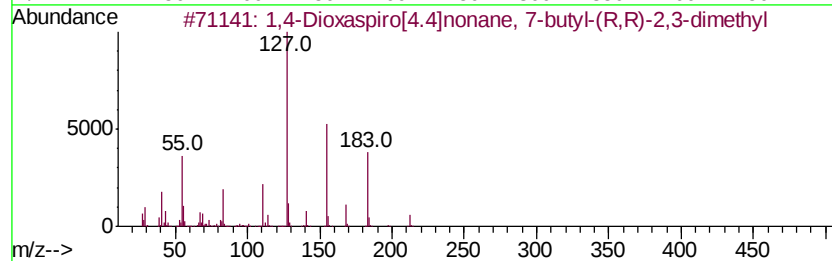
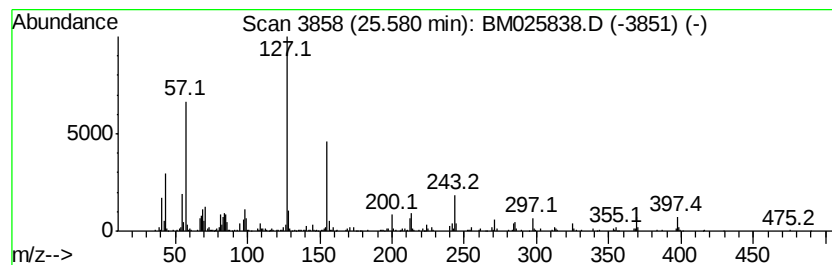
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 1,4-Dioxaspiro[4.4]nonane, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.58	7.58 ng/ul	1273240	Perylene-d12	23.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,4-Dioxaspiro[4.4]nonane, 7-but...	212	C13H24O2	104807-77-4	64
2		Diethylmalonic acid, 3,4-difluor...	328	C17H22F2O4	1000369-32-3	38
3		Urea, 1-(4-chlorophenyl)-3-propyl-	212	C10H13ClN2O	013208-64-5	38
4		2-Hydroxy-4-hydroxylaminopyrimidine	127	C4H5N3O2	020555-88-8	38
5		1H-Imidazole, 1-methyl-4-nitro-	127	C4H5N3O2	003034-41-1	38



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032620\
Data File : BM025838.D
Acq On : 27 Mar 2020 15:15
Operator : CG/JU
Sample : L2005-08
Misc :
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Instrument :
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ClientSampleId :
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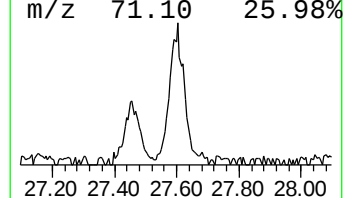
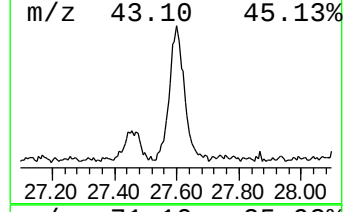
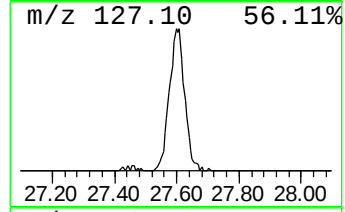
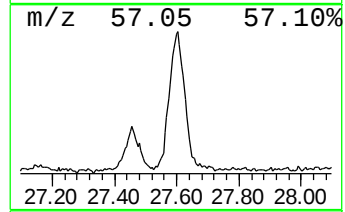
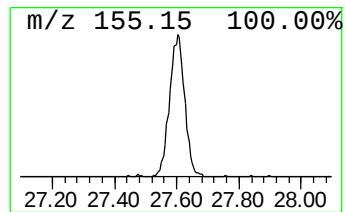
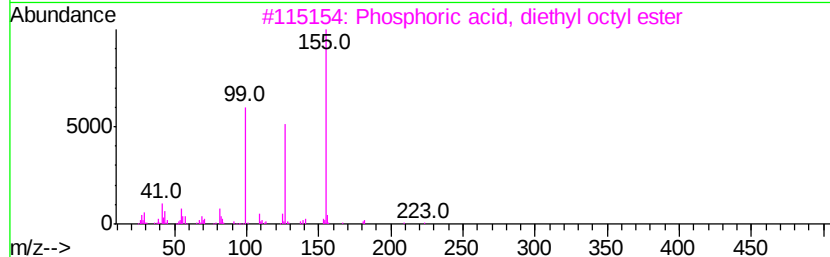
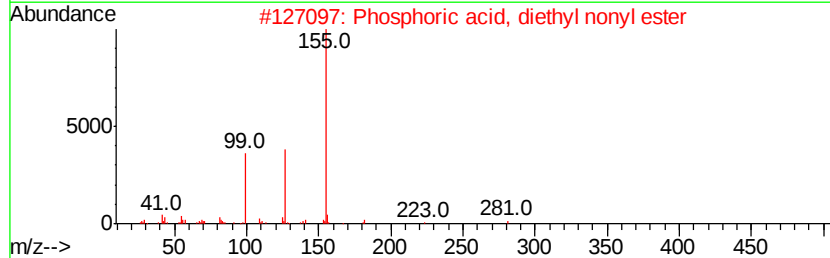
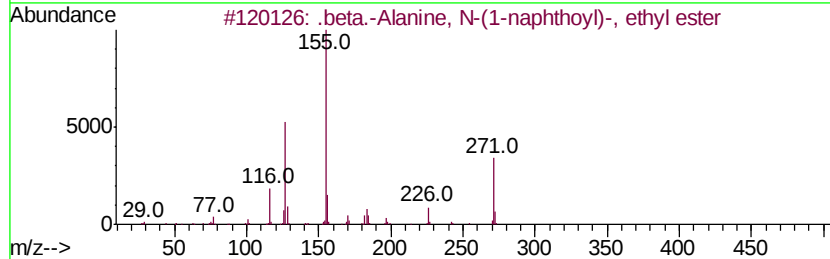
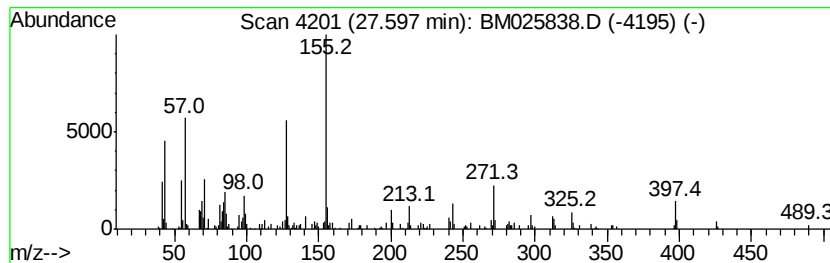
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 .beta.-Alanine, N-(1-naphth... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.60	3.83 ng/ul	642745	Perylene-d12	23.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	.beta.-Alanine, N-(1-naphthovl)-...	271	C16H17N03	1000321-94-4	72
2		Phosphoric acid, diethyl nonyl e...	280	C13H29O4P	1000308-95-2	53
3		Phosphoric acid, diethyl octyl e...	266	C12H27O4P	1000308-90-4	53
4		Imidazole, 4-amino-5-ethoxycarbo...	155	C6H9N3O2	021190-16-9	53
5		5-Oxazolidinone, 2-(1,1-dimethyl...	297	C18H19N03	119215-56-4	46



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032620\
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 Sample : L2005-08
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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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	10.1	ng/ul	783938	2	10.35	1550810	20.0
o-Chloroaniline	9.38	76.1	ng/ul	5902130	2	10.35	1550810	20.0
Phenol, 2,5-dichl...	10.03	8.5	ng/ul	656244	2	10.35	1550810	20.0
Phenol, 2,3-dichl...	10.13	15.8	ng/ul	1222190	2	10.35	1550810	20.0
Benzene, 1-chloro...	10.93	12.0	ng/ul	927450	2	10.35	1550810	20.0
2-Butenoic acid, ...	12.00	8.3	ng/ul	643223	2	10.35	1550810	20.0
Benzenamine, 2,5-...	12.62	6.2	ng/ul	637115	3	14.20	2044100	20.0
L-Valine, N-dimet...	24.08	4.2	ng/ul	700306	6	23.29	3360470	20.0
1,4-Dioxaspiro[4...	25.58	7.6	ng/ul	1273240	6	23.29	3360470	20.0
.beta.-Alanine, N...	27.60	3.8	ng/ul	642745	6	23.29	3360470	20.0