

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
 Data File : BM004645.D
 Acq On : 16 Mar 2016 21:07
 Operator : UM/SJ
 Sample : H1783-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB2

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 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:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.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	2.993	46	52	65	rBV	1026637	1487832	2.49%	0.645%
2	5.293	417	443	447	rBV	8422067	59672052	100.00%	25.881%
3	7.775	851	865	875	rBV2	6100518	23082760	38.68%	10.012%
4	7.969	875	898	902	rBV	7057252	38421766	64.39%	16.665%
5	8.298	931	954	959	rBV	6925056	40641772	68.11%	17.628%
6	9.345	1124	1132	1141	rBV	1430325	2477471	4.15%	1.075%
7	9.569	1163	1170	1174	rBV	410076	704131	1.18%	0.305%
8	9.616	1174	1178	1180	rVV	373046	629463	1.05%	0.273%
9	9.657	1180	1185	1191	rVV	810458	1534688	2.57%	0.666%
10	10.287	1279	1292	1296	rBV2	392762	1290654	2.16%	0.560%
11	10.545	1322	1336	1337	rBV2	6115077	20265261	33.96%	8.790%
12	11.063	1411	1424	1430	rBV3	6065578	20113298	33.71%	8.724%
13	12.633	1682	1691	1693	rBV	352200	612618	1.03%	0.266%
14	12.669	1693	1697	1704	rVB	1033605	1614231	2.71%	0.700%
15	13.280	1793	1801	1812	rVB	3681695	6494423	10.88%	2.817%
16	13.886	1894	1904	1909	rBV	619590	1083885	1.82%	0.470%
17	14.169	1946	1952	1959	rBV	712729	1052682	1.76%	0.457%
18	14.480	1998	2005	2013	rBV2	394438	697125	1.17%	0.302%
19	15.469	2166	2173	2186	rVB	880246	1340649	2.25%	0.581%
20	17.221	2465	2471	2476	rBV	470964	677523	1.14%	0.294%
21	17.321	2481	2488	2498	rBV	957277	1386191	2.32%	0.601%
22	19.609	2871	2877	2888	rBV	1203140	1686513	2.83%	0.731%
23	21.403	3176	3182	3188	rBV	682569	899140	1.51%	0.390%
24	23.556	3540	3548	3560	rBV	926283	1816193	3.04%	0.788%
25	23.703	3566	3573	3584	rVB2	443880	876565	1.47%	0.380%

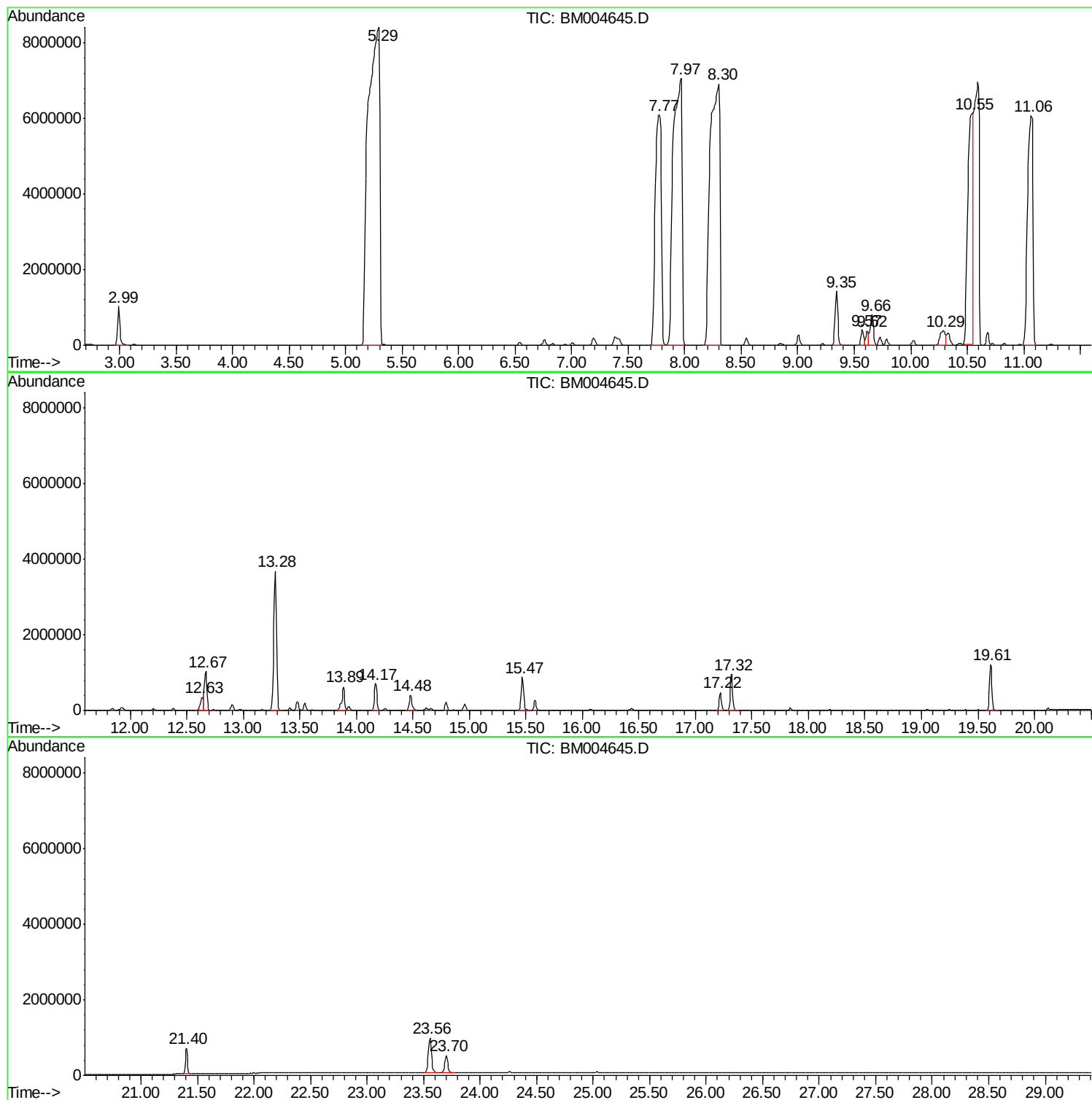
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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
Quant Title : SVOA CALIBRATION

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



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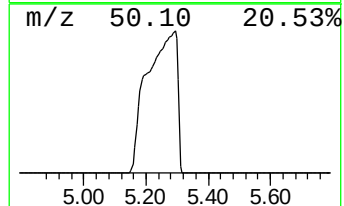
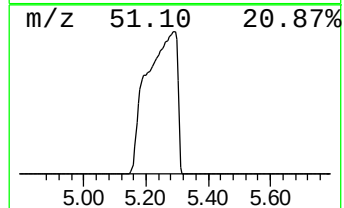
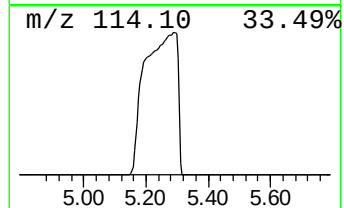
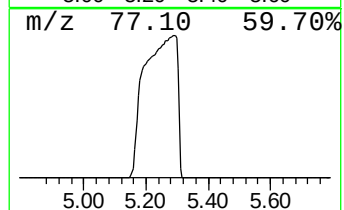
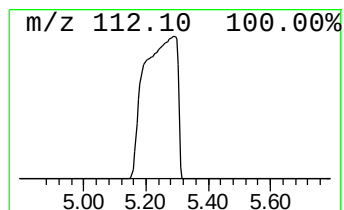
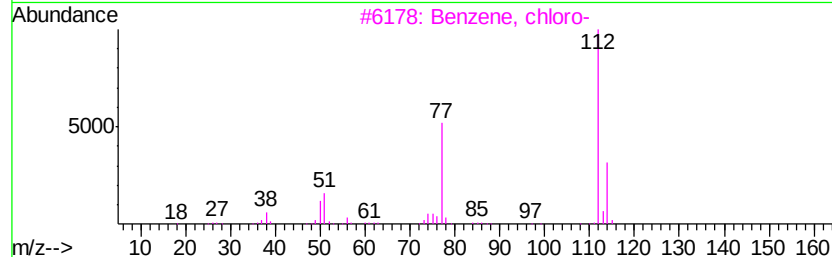
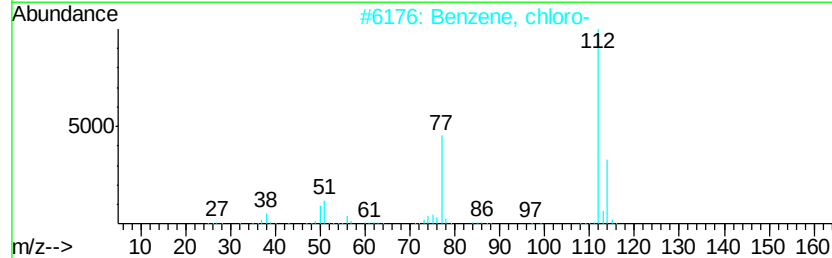
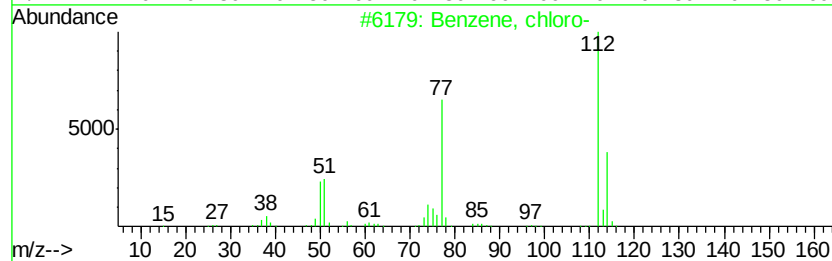
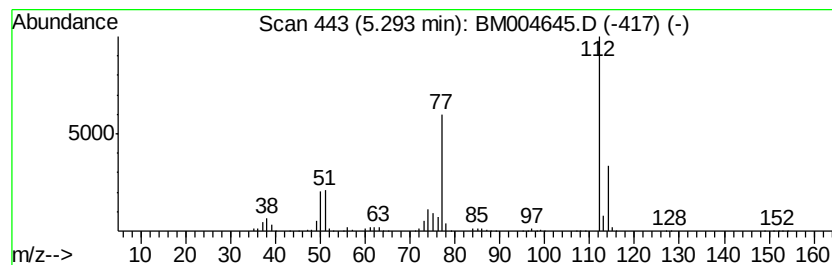
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.29	51.70 ng/ul	59672100	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	96
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	91
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	90
5		Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	27



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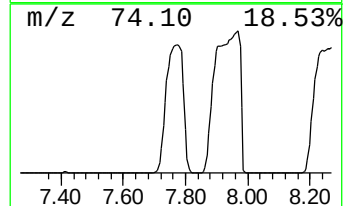
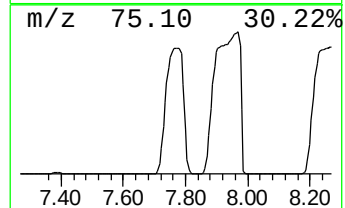
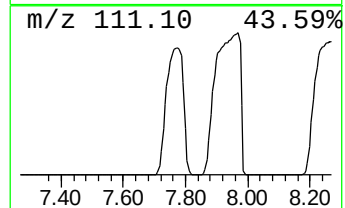
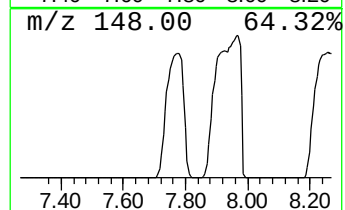
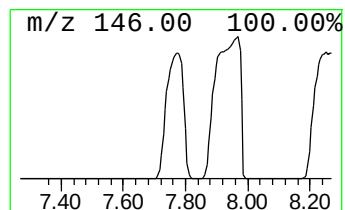
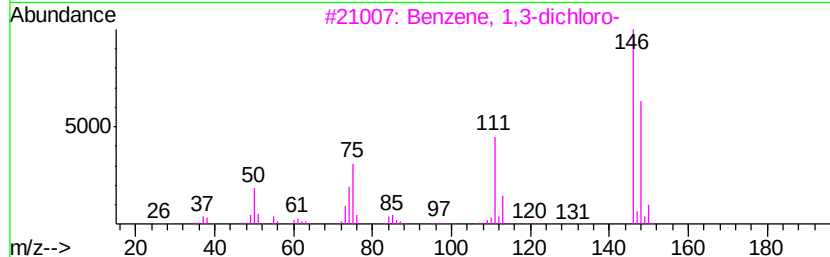
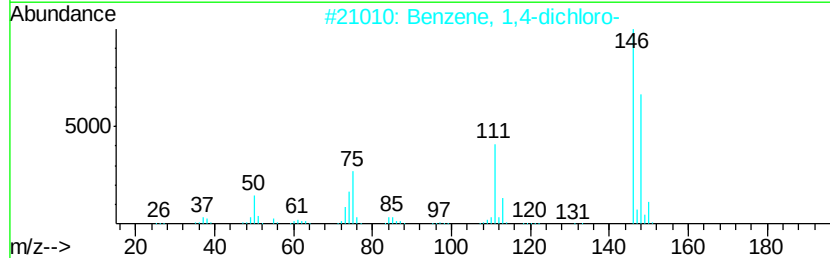
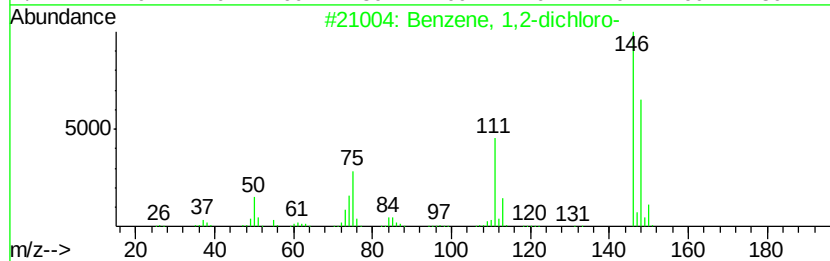
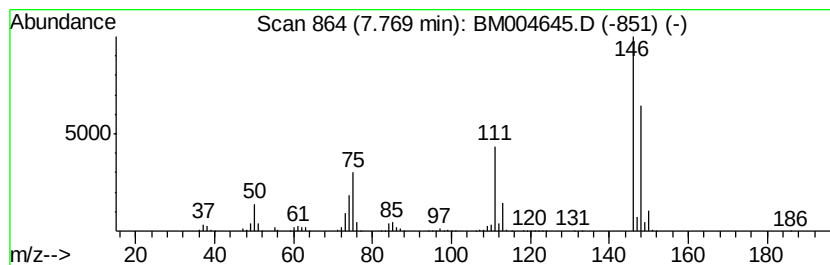
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.77	20.00 ng/ul	23082800	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96



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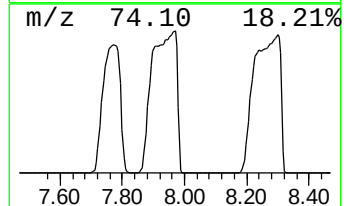
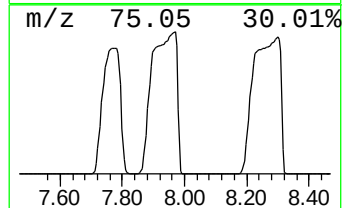
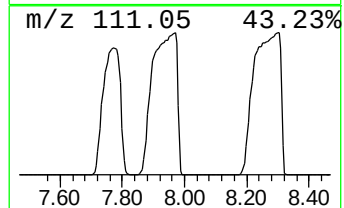
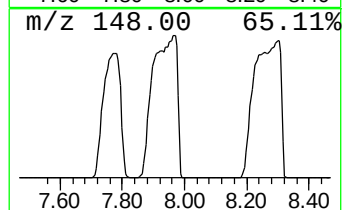
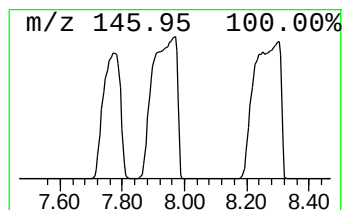
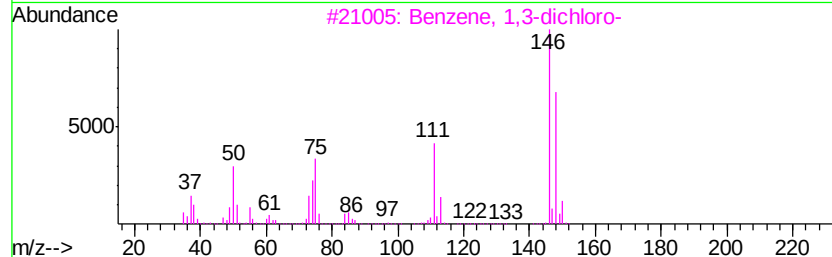
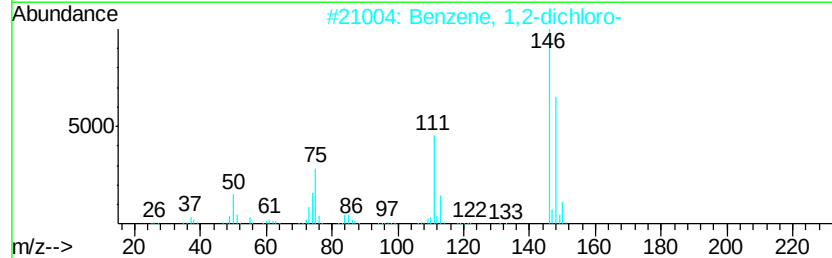
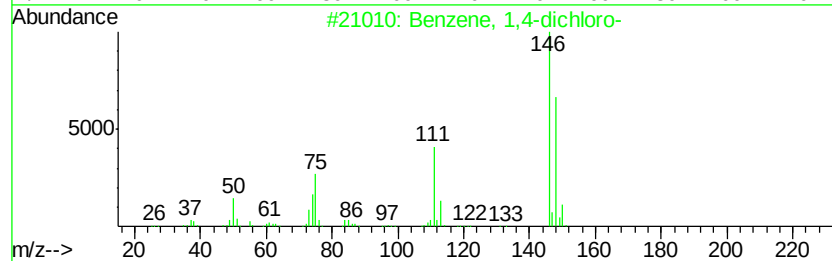
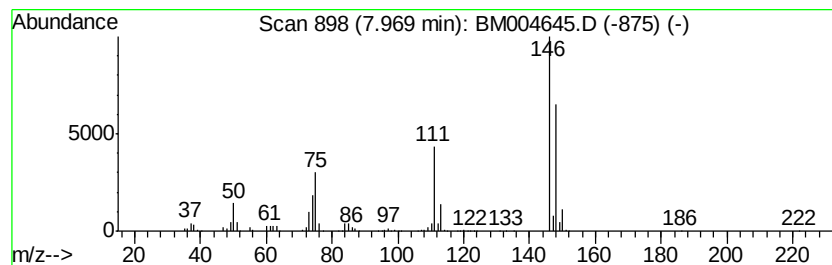
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.97	33.29 ng/ul	38421800	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	97
5		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	97



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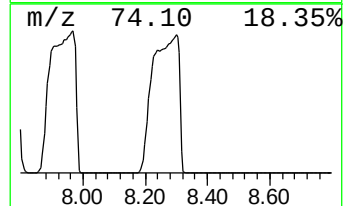
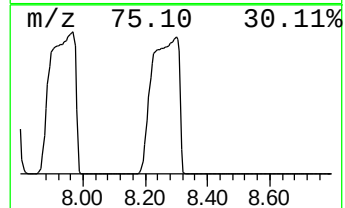
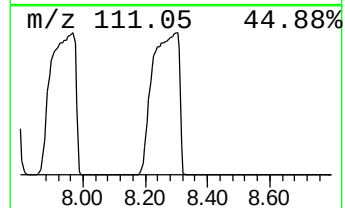
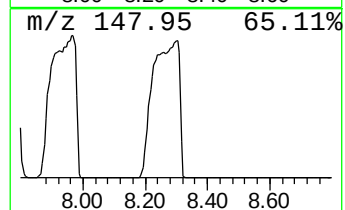
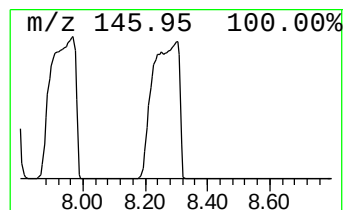
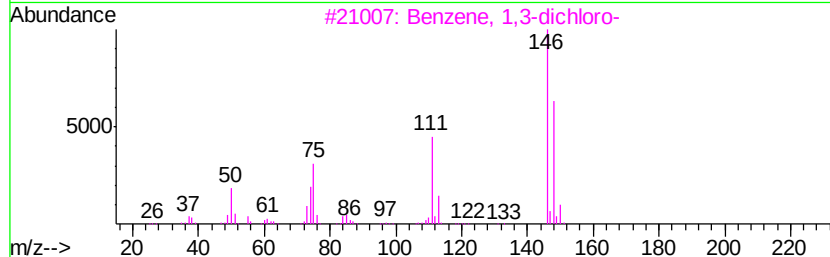
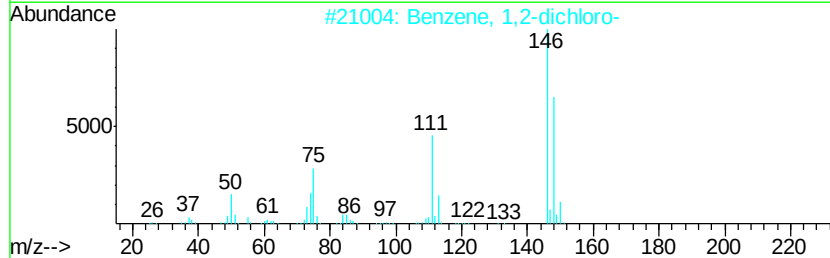
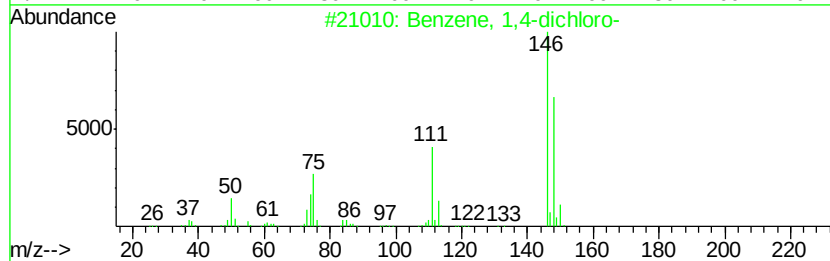
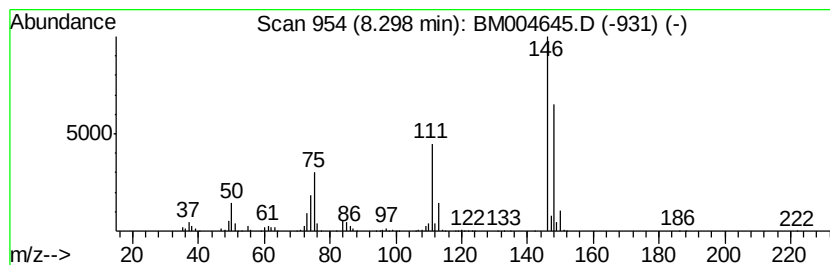
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Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,3-dichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.30	35.21 ng/ul	40641800	1,4-Dichlorobenzene-d4	7.87

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	98
2		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	98
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
4		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96



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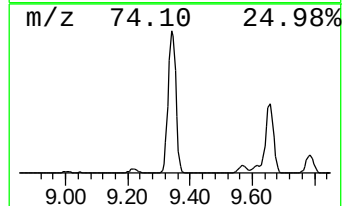
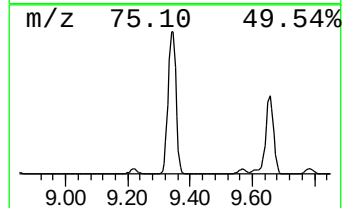
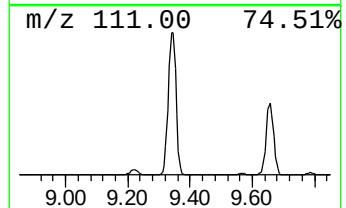
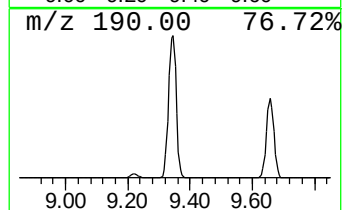
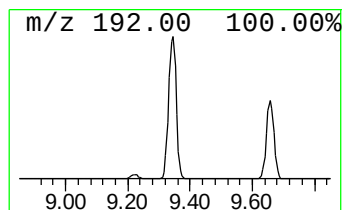
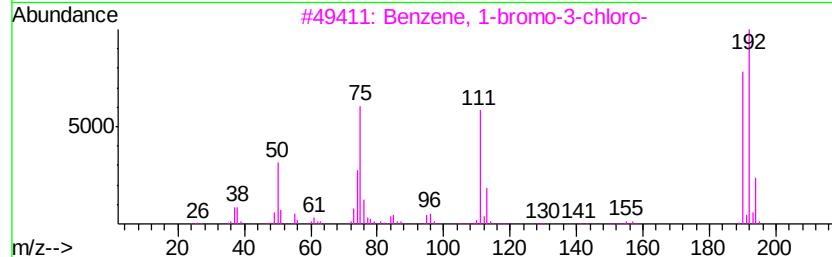
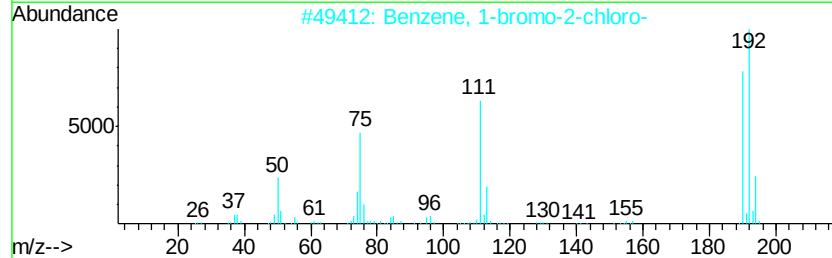
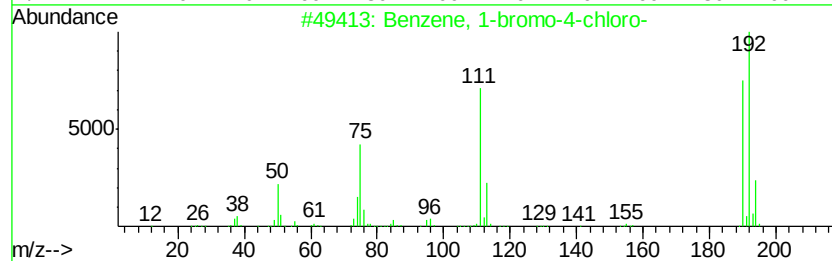
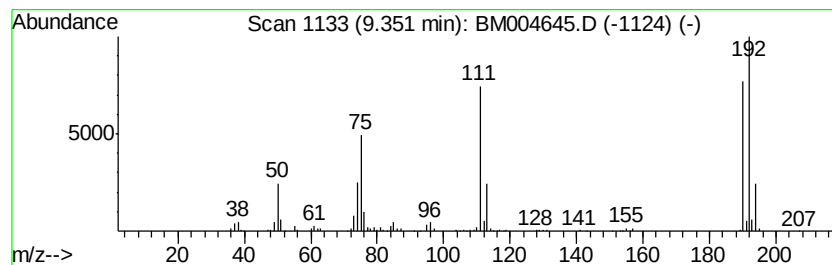
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Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.35	2.45 ng/ul	2477470	Napthalene-d8	10.68

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	96
5		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	96



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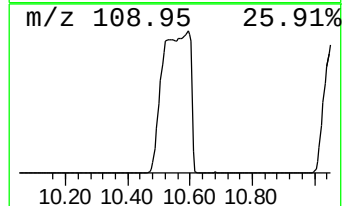
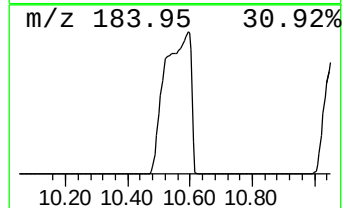
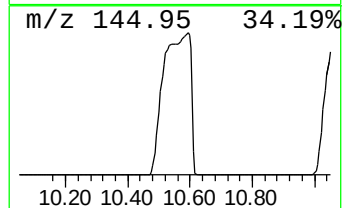
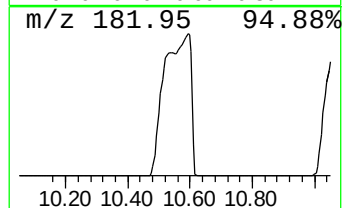
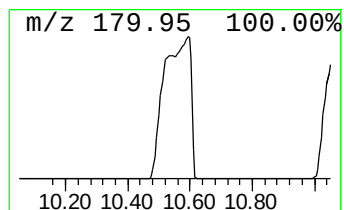
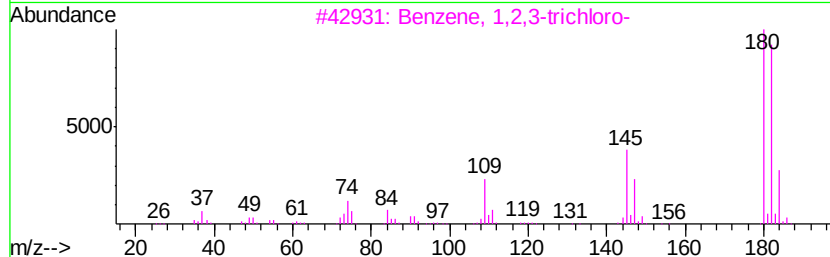
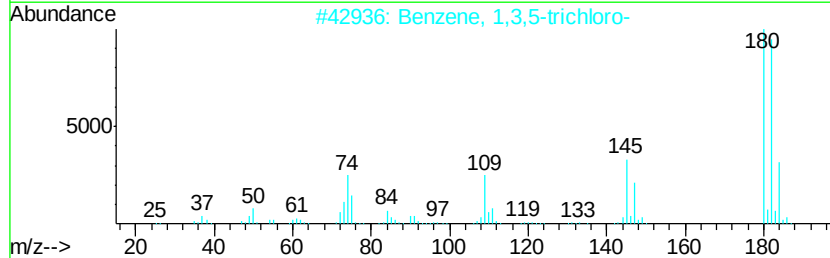
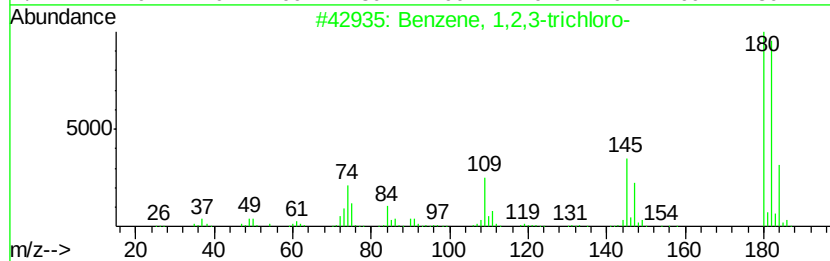
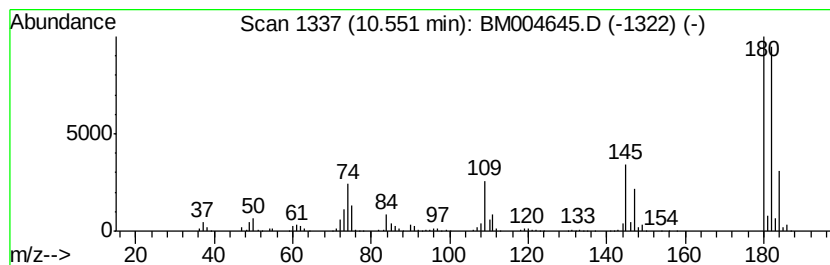
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BNA_M
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,2,3-trichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.55	20.00 ng/ul	20265300	Naphthalene-d8	10.68
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6 99
2		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 98
3		Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6 98
4		Benzene, 1,2,4-trichloro-	180 C6H3Cl3	000120-82-1 97
5		Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6 97



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
Data File : BM004645.D
Acq On : 16 Mar 2016 21:07
Operator : UM/SJ
Sample : H1783-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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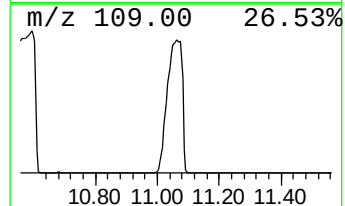
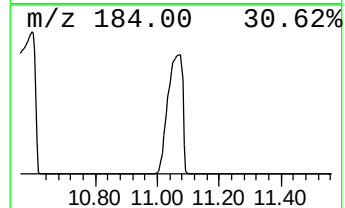
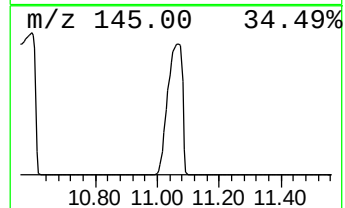
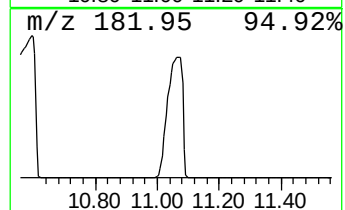
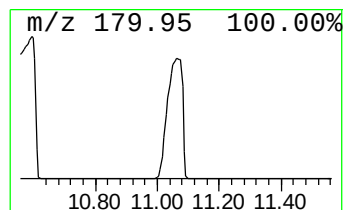
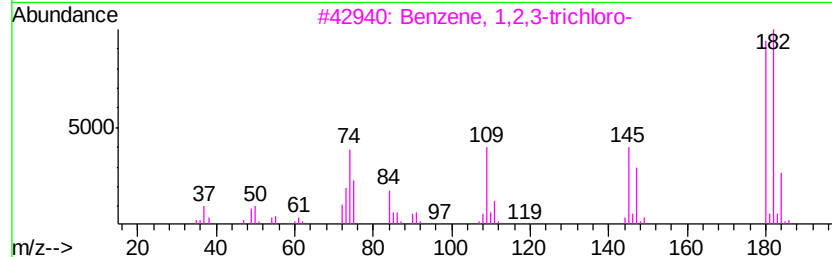
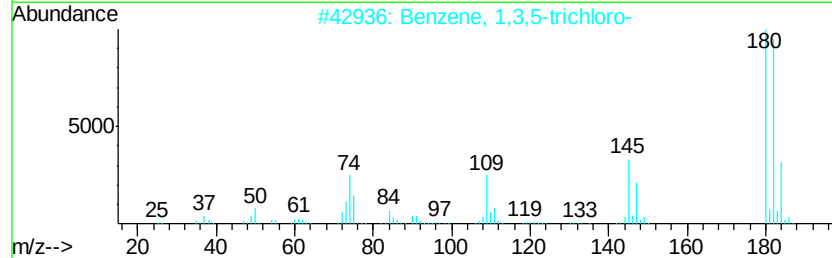
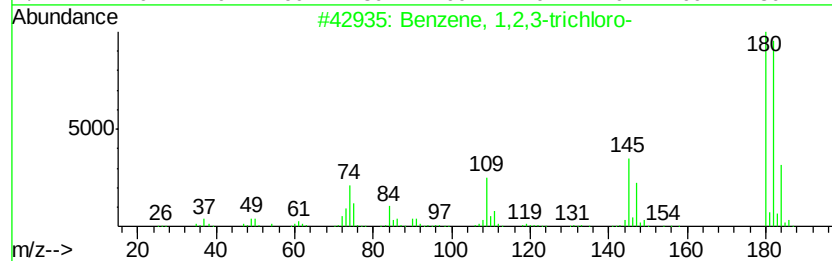
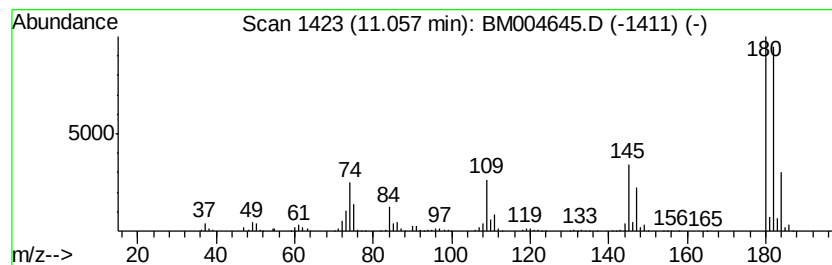
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,3,5-trichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.06	19.85 ng/ul	20113300	Napthalene-d8	10.68

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	99
2		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3		Benzene, 1,2,3-trichloro-	180	C6H3Cl3	000087-61-6	97
4		Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5		Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	96



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
Data File : BM004645.D
Acq On : 16 Mar 2016 21:07
Operator : UM/SJ
Sample : H1783-11
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AB2

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, chloro-	5.29	51.7	ng/ul	59672100	1	7.87	23082800	20.0
Benzene, 1,2-dich...	7.77	20.0	ng/ul	23082800	1	7.87	23082800	20.0
Benzene, 1,4-dich...	7.97	33.3	ng/ul	38421800	1	7.87	23082800	20.0
Benzene, 1,3-dich...	8.30	35.2	ng/ul	40641800	1	7.87	23082800	20.0
Benzene, 1-bromo-...	9.35	2.5	ng/ul	2477470	2	10.68	20265300	20.0
Benzene, 1,2,3-tr...	10.55	20.0	ng/ul	20265300	2	10.68	20265300	20.0
Benzene, 1,3,5-tr...	11.06	19.9	ng/ul	20113300	2	10.68	20265300	20.0