

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
 Data File : BM019375.D
 Acq On : 24 Mar 2019 05:00
 Operator : JU/SJ
 Sample : K2025-14
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0985

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-BM032219MA.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	4.969	327	337	351	rBV2	40958093	73877298	63.97%	17.325%
2	6.793	642	647	661	rBV2	118001	242893	0.21%	0.057%
3	6.963	670	676	691	rBV	449041	743603	0.64%	0.174%
4	7.146	701	707	711	rBV	494978	831702	0.72%	0.195%
5	7.499	759	767	780	rBV	14817547	24027620	20.80%	5.635%
6	7.669	780	796	801	rBV2	43643579	97190225	84.15%	22.792%
7	7.993	837	851	856	rBV2	45063364	115490808	100.00%	27.084%
8	8.322	902	907	921	rBV	318020	577114	0.50%	0.135%
9	8.787	978	986	987	rBV	452435	703120	0.61%	0.165%
10	8.846	987	996	1016	rVB	13210966	26802973	23.21%	6.286%
11	9.099	1033	1039	1052	rVB2	125687	221085	0.19%	0.052%
12	9.493	1101	1106	1119	rVB	423555	777600	0.67%	0.182%
13	10.022	1189	1196	1203	rBV2	657859	1395791	1.21%	0.327%
14	10.087	1203	1207	1217	rVB	479315	842239	0.73%	0.198%
15	10.181	1217	1223	1228	rBV	268400	456314	0.40%	0.107%
16	10.263	1228	1237	1246	rBV	17513466	28154450	24.38%	6.603%
17	10.393	1253	1259	1266	rVB	868005	1393199	1.21%	0.327%
18	10.528	1276	1282	1288	rBV2	64038	158709	0.14%	0.037%
19	10.775	1315	1324	1336	rBV	7796783	12237709	10.60%	2.870%
20	11.010	1356	1364	1372	rBV	4915201	8052446	6.97%	1.888%
21	11.222	1391	1400	1419	rBV	1084724	2139386	1.85%	0.502%
22	12.434	1595	1606	1611	rBV	148989	309348	0.27%	0.073%
23	12.851	1671	1677	1683	rBV	102928	169290	0.15%	0.040%
24	13.051	1705	1711	1717	rBV	587786	944604	0.82%	0.222%
25	13.298	1747	1753	1765	rBV	318103	604654	0.52%	0.142%
26	13.687	1812	1819	1826	rBV	1196584	1705370	1.48%	0.400%
27	13.957	1858	1865	1877	rBV	1424793	2083849	1.80%	0.489%
28	14.269	1909	1918	1926	rBV2	1175718	1829493	1.58%	0.429%
29	15.263	2081	2087	2104	rBV	1767063	2571141	2.23%	0.603%
30	15.404	2106	2111	2125	rBV	533205	839749	0.73%	0.197%
31	17.022	2380	2386	2396	rBV2	1555679	2117862	1.83%	0.497%
32	17.122	2396	2403	2417	rBV	2044142	2907200	2.52%	0.682%
33	19.427	2789	2795	2806	rBV	2645793	3541921	3.07%	0.831%
34	21.227	3096	3101	3111	rBV	1920323	2629321	2.28%	0.617%

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Title : SVOA CALIBRATION

35	22.245	3271	3274	3280	rVB	115801	140139	0.12%	0.033%
36	22.739	3355	3358	3363	rVB	143629	190671	0.17%	0.045%
37	23.298	3447	3453	3469	rVB2	2304424	4240763	3.67%	0.995%
38	23.445	3471	3478	3491	rVV	1443323	2766799	2.40%	0.649%
39	23.927	3556	3560	3566	rVB	151979	268427	0.23%	0.063%
40	24.656	3681	3684	3690	rVB	142243	242062	0.21%	0.057%

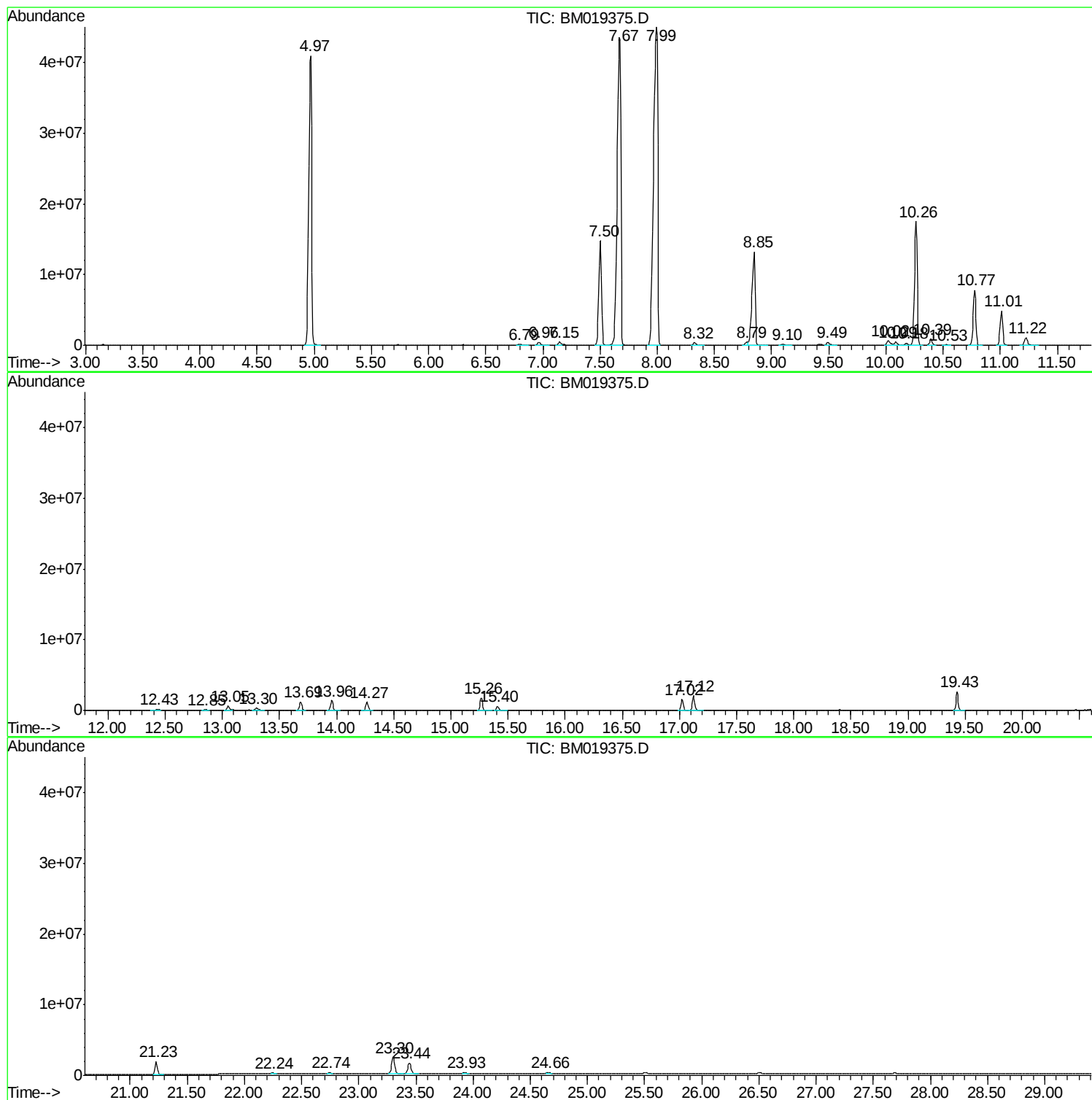
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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



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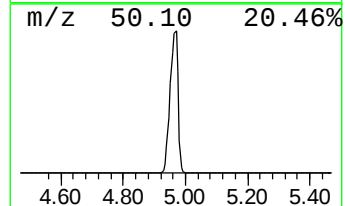
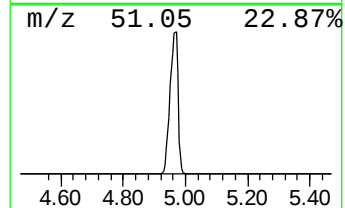
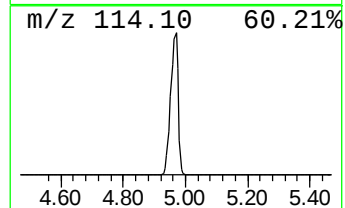
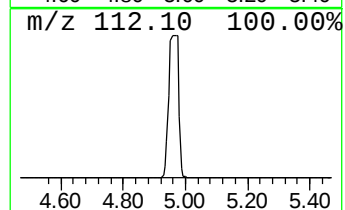
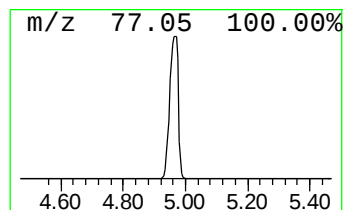
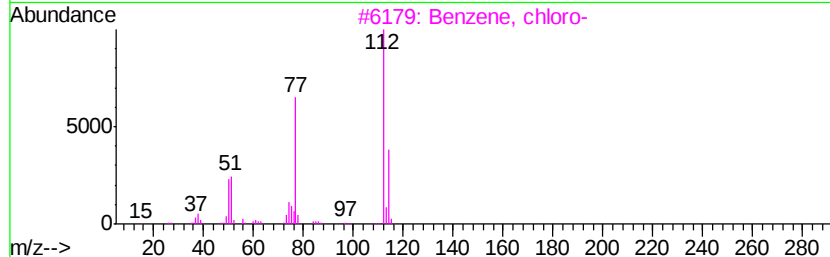
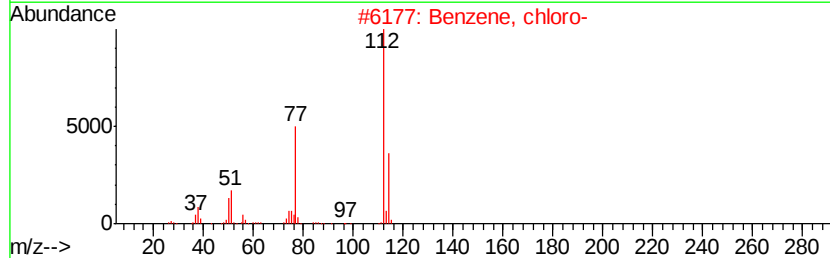
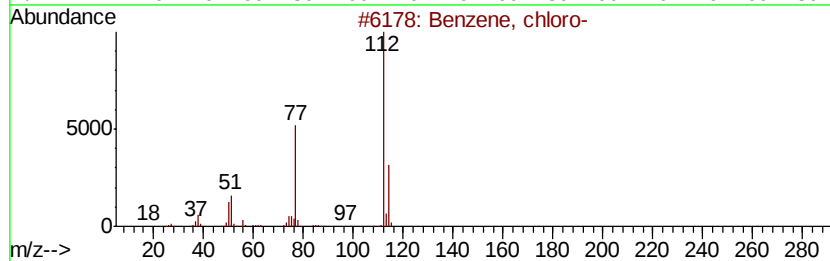
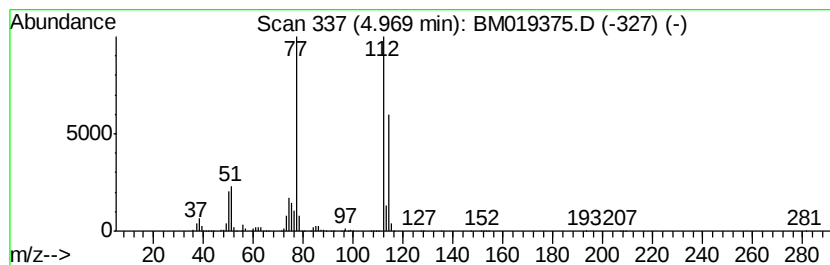
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	15.20 ng/ul	73877300	1,4-Dichlorobenzene-d4	7.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
2			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
3			Benzene, chloro-	112	C6H5Cl	000108-90-7	90
4			Benzene, chloro-	112	C6H5Cl	000108-90-7	86
5			Phenol, 3-fluoro-	112	C6H5FO	000372-20-3	14



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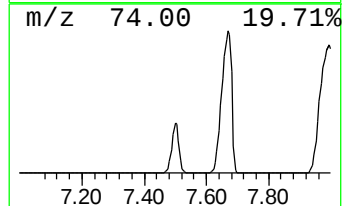
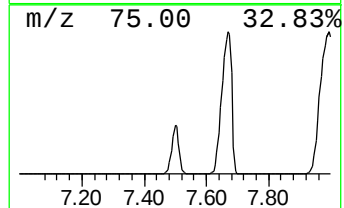
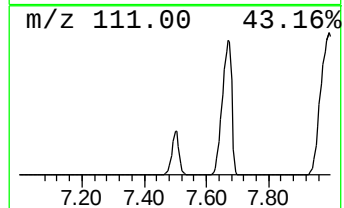
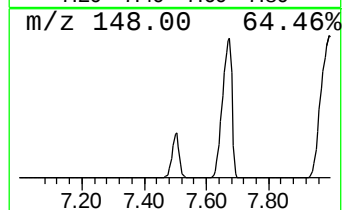
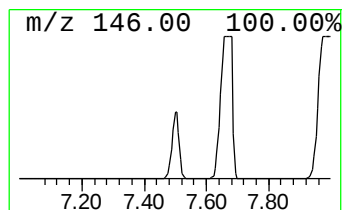
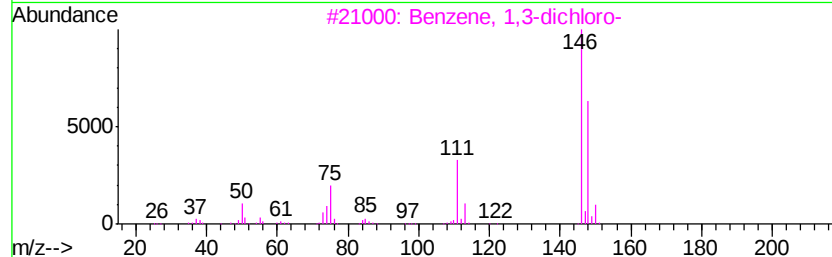
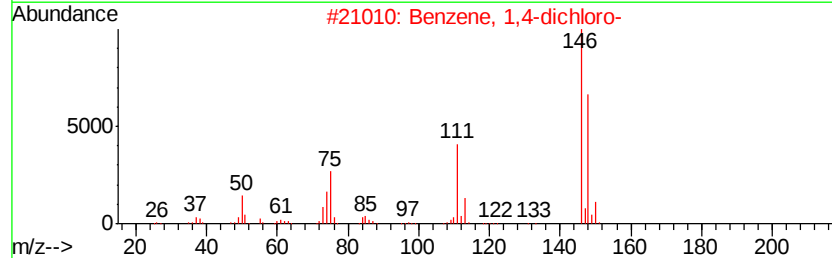
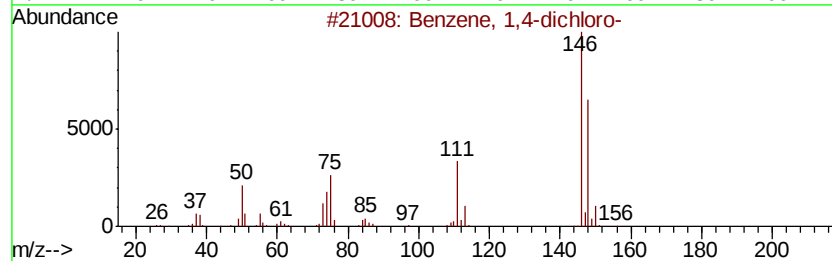
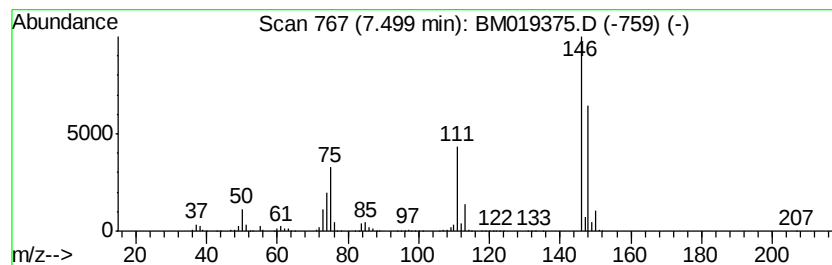
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.50	4.94 ng/ul	24027600	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95
4		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	95



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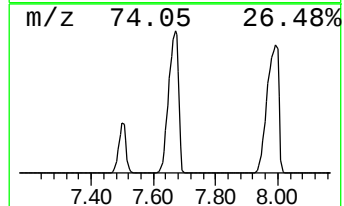
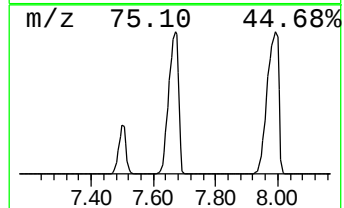
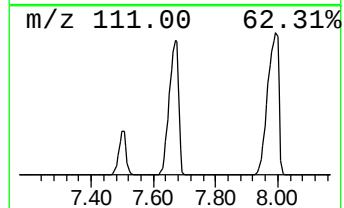
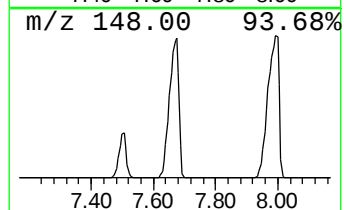
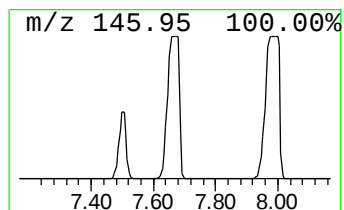
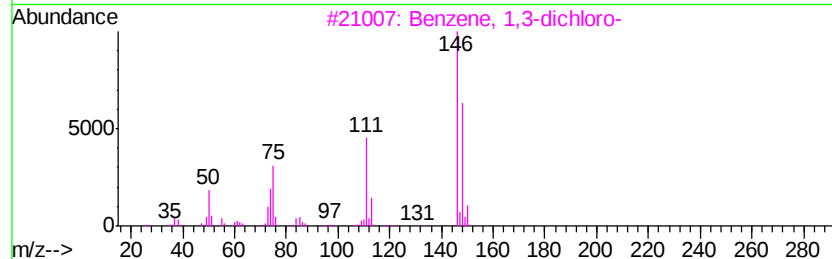
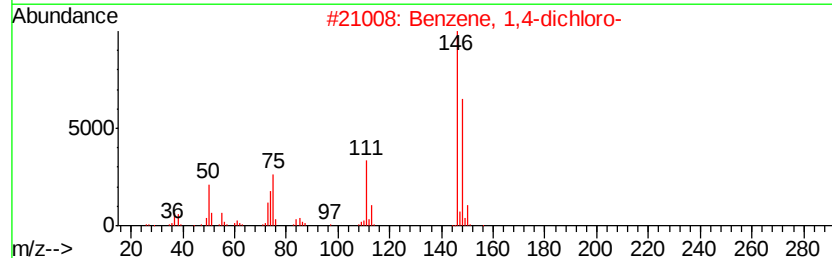
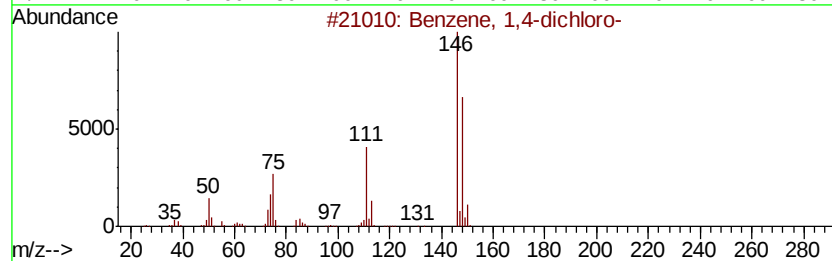
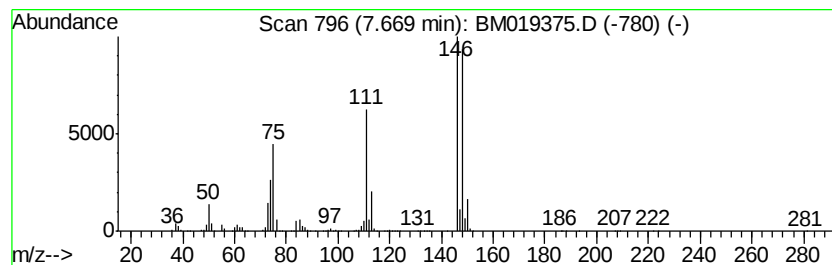
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	20.00 ng/ul	97190200	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	91
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	91



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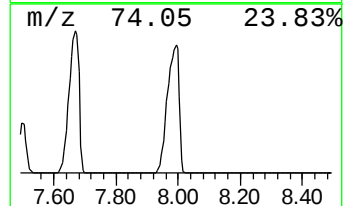
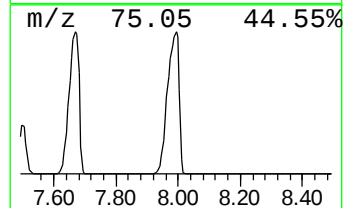
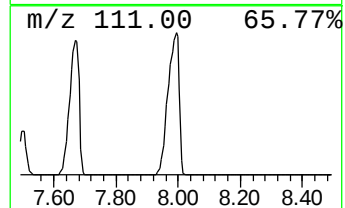
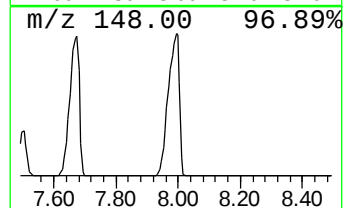
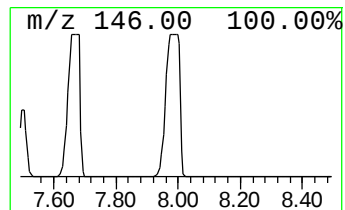
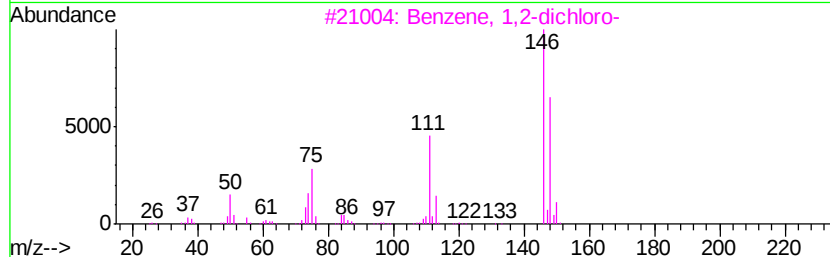
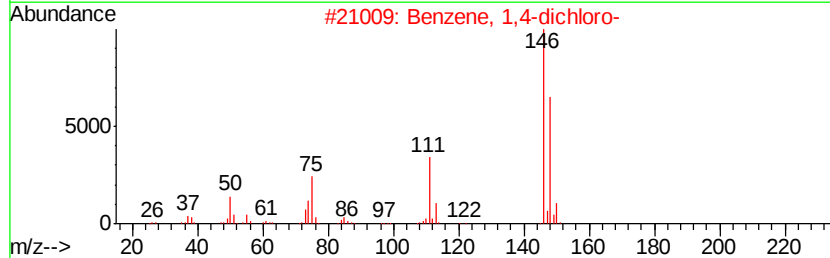
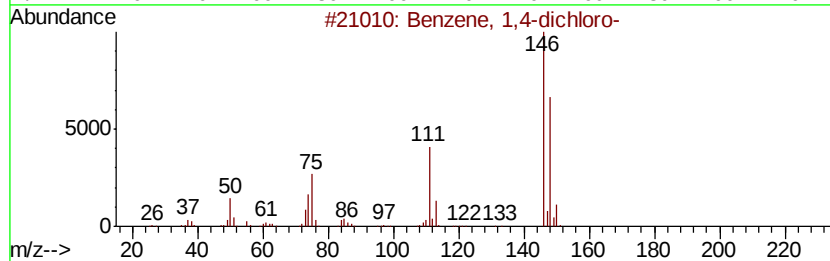
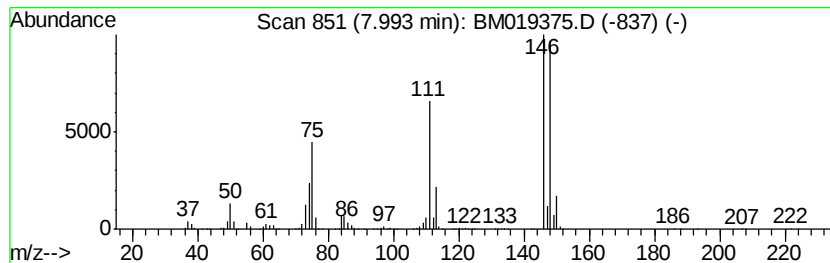
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Peak Number 4 Benzene, 1,2-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.99	23.77 ng/ul	115491000	1,4-Dichlorobenzene-d4	7.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	94
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	87



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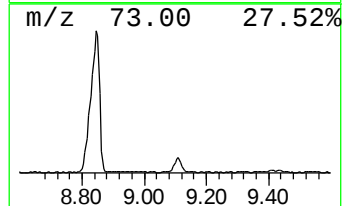
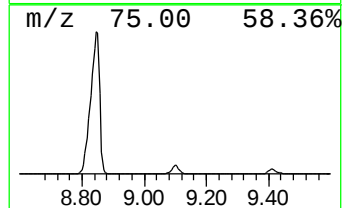
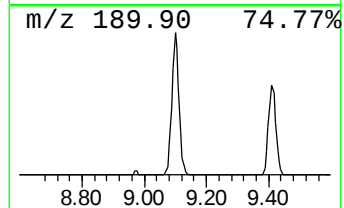
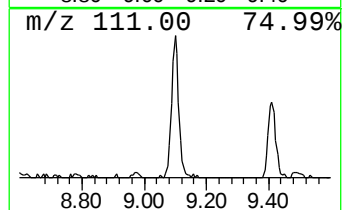
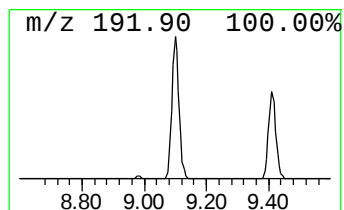
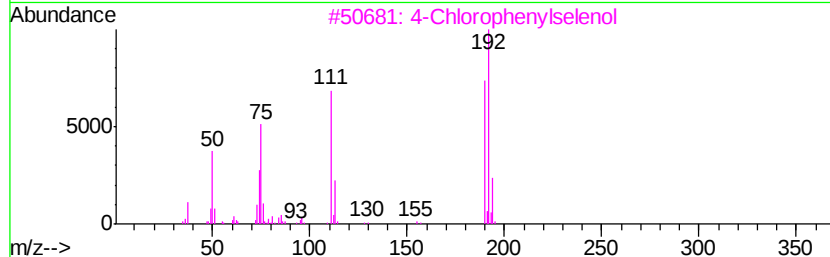
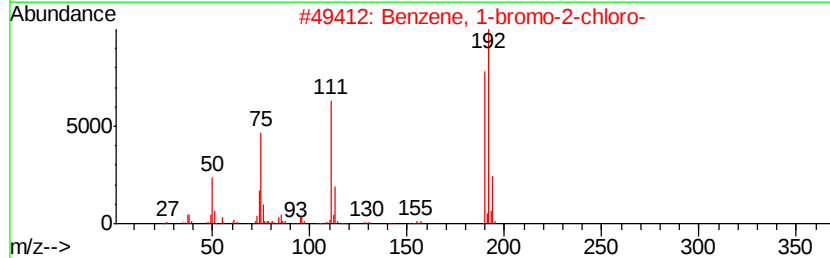
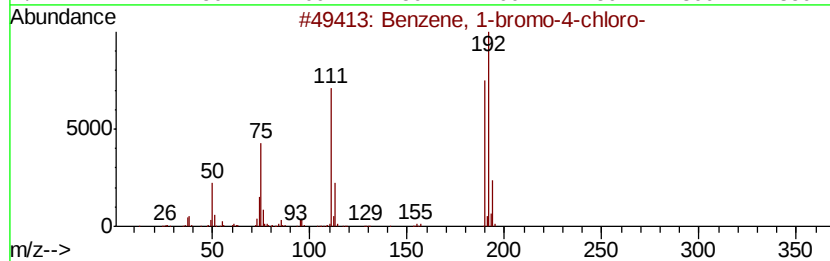
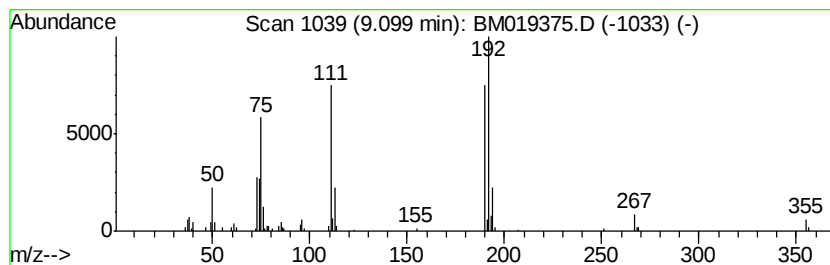
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-bromo-4-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.10	3.17 ng/ul	221085	Napthalene-d8	10.39

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		4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
4		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	96
5		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
Sample : K2025-14
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0985

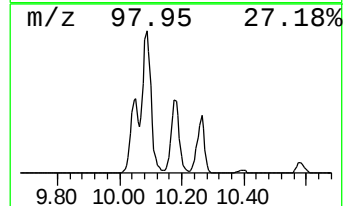
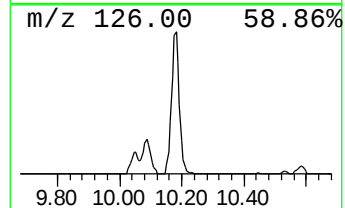
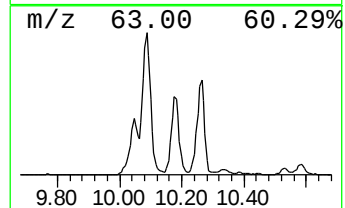
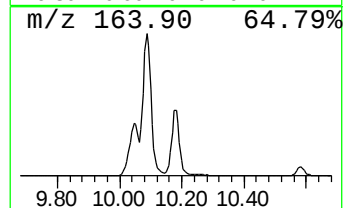
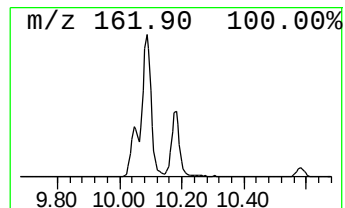
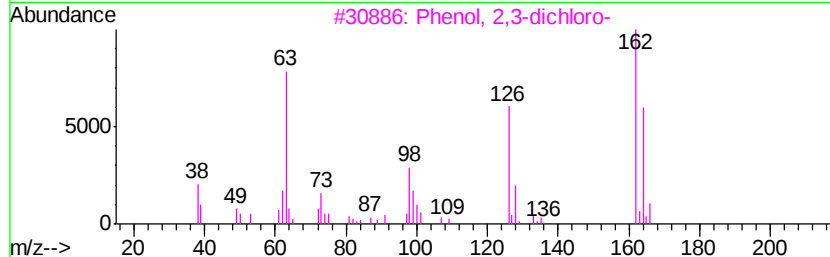
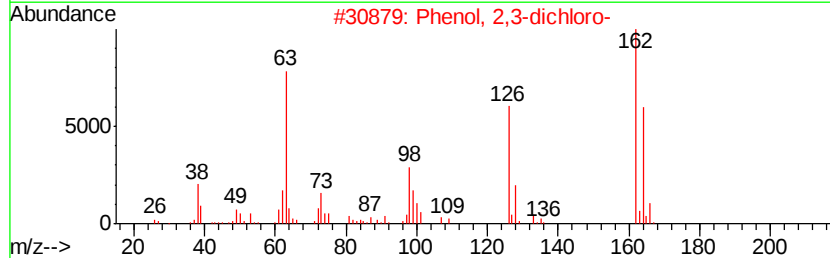
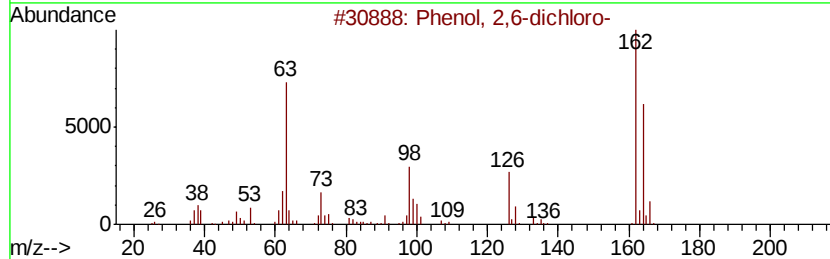
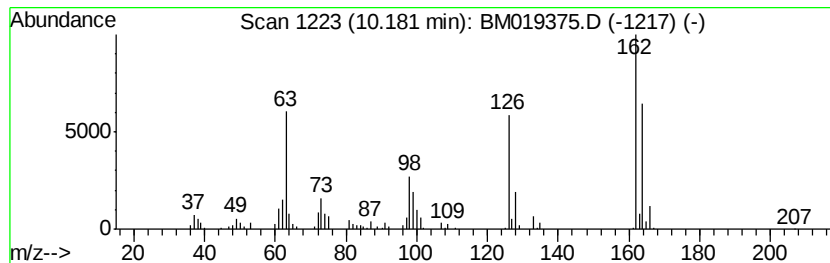
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 2,6-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.18	6.55 ng/ul	456314	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	96
2		Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	95
3		Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	95
4		Phenol, 2,3-dichloro-	162	C6H4Cl2O	000576-24-9	93
5		Phenol, 2,6-dichloro-	162	C6H4Cl2O	000087-65-0	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
Sample : K2025-14
Misc :
ALS Vial : 33 Sample Multiplier: 1

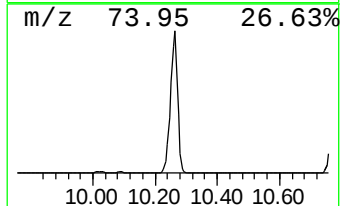
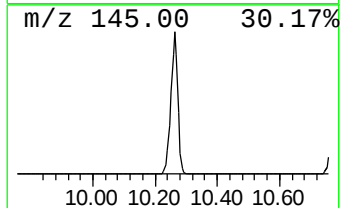
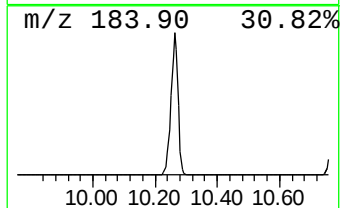
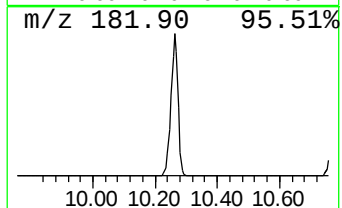
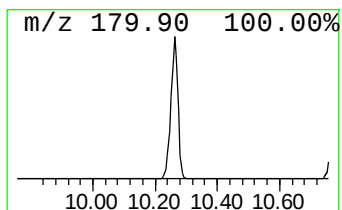
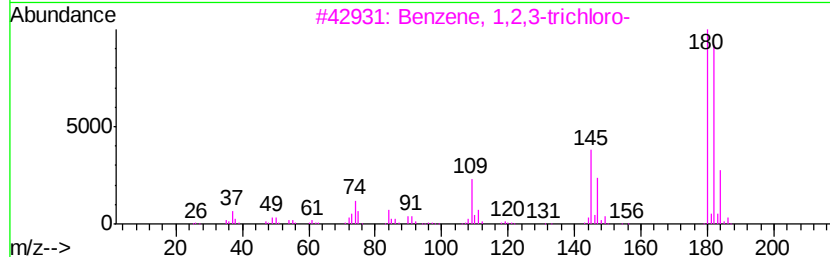
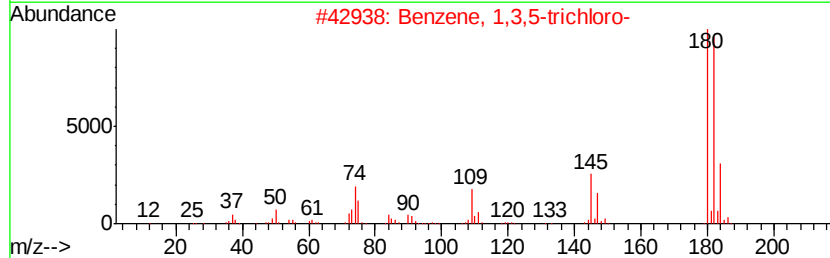
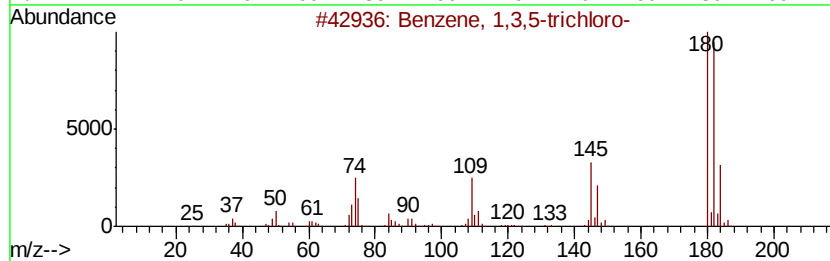
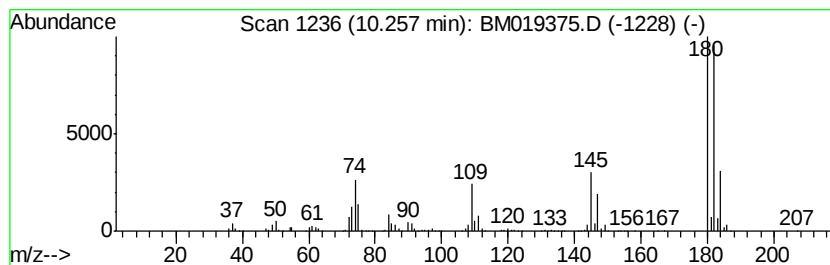
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.26	404.17 ng/ul	28154500	Napthalene-d8	10.39
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 98
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, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3 96
5		Benzene, 1,2,4-trichloro-	180 C6H3Cl3	000120-82-1 95



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
Sample : K2025-14
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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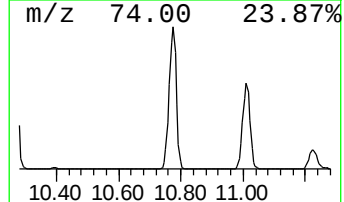
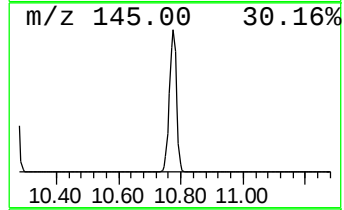
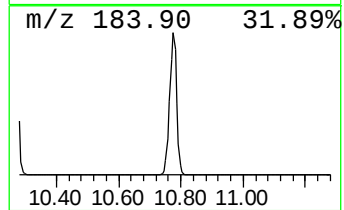
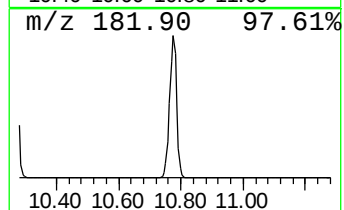
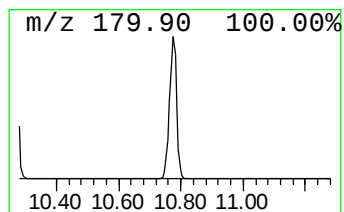
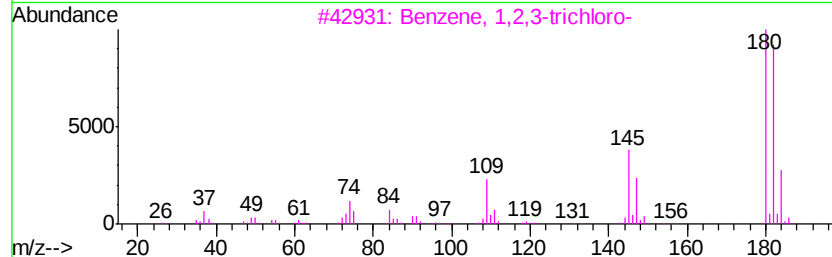
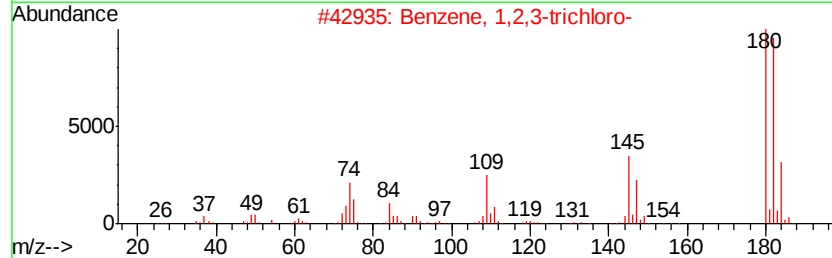
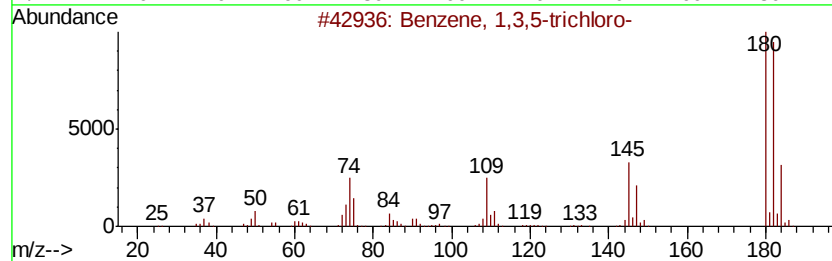
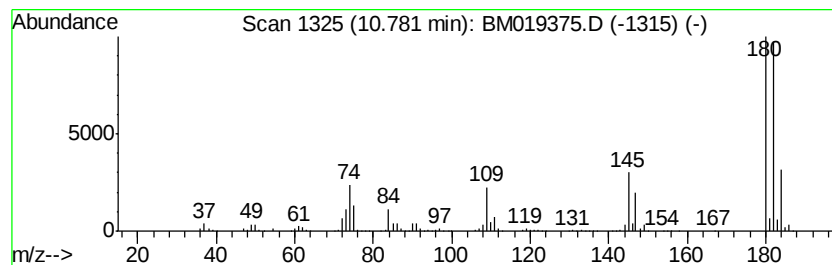
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,3-trichloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	175.68 ng/ul	12237700	Naphthalene-d8	10.39

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
Sample : K2025-14
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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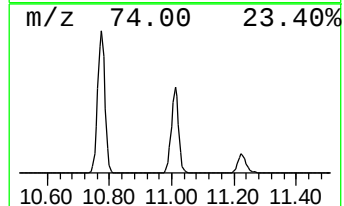
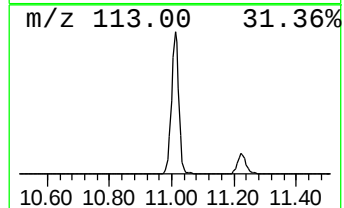
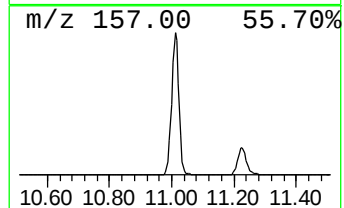
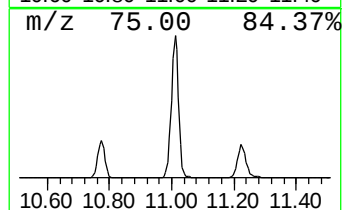
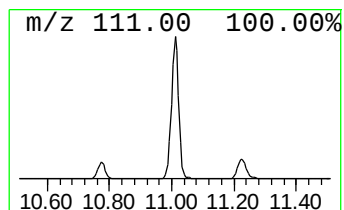
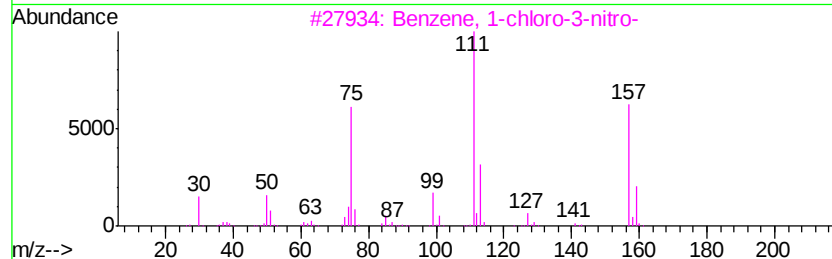
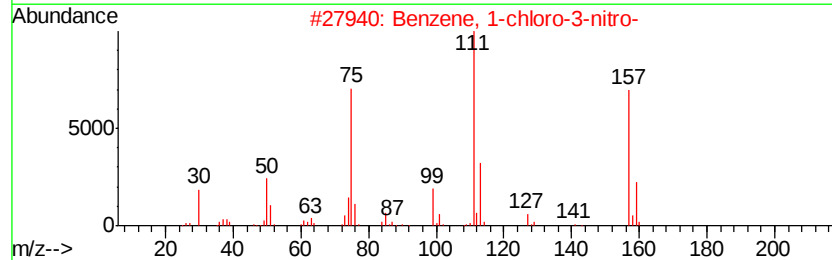
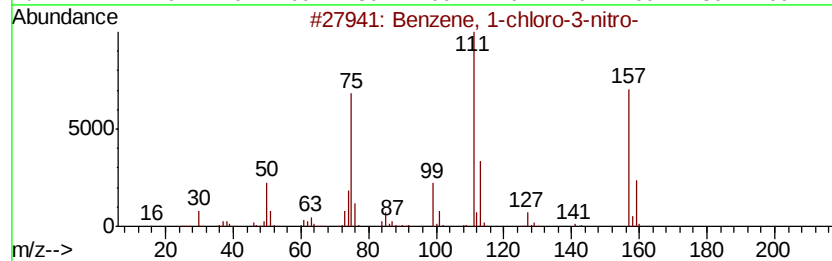
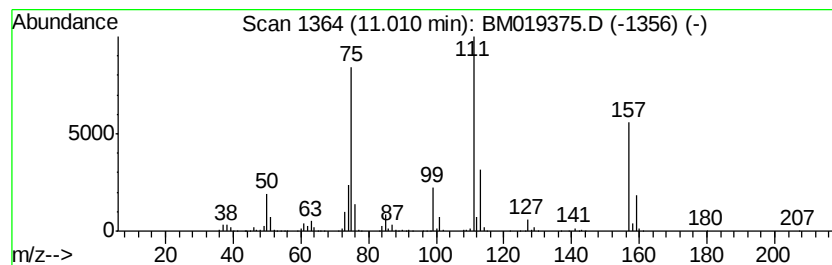
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.01	115.60 ng/ul	8052450	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	95
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	83



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
Sample : K2025-14
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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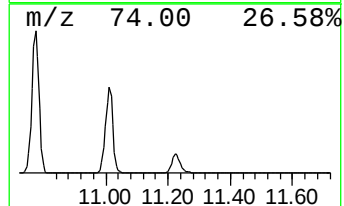
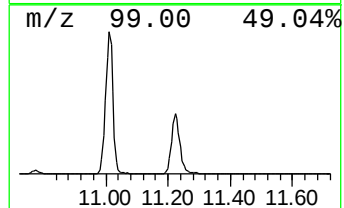
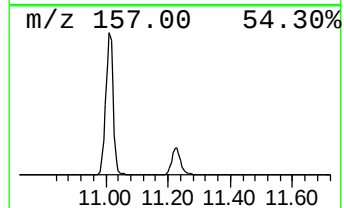
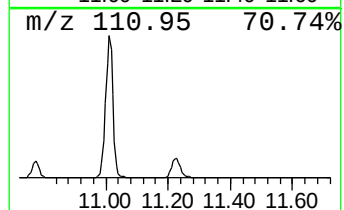
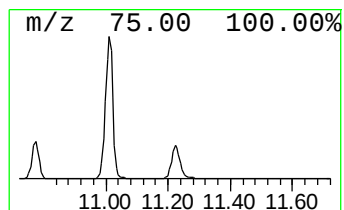
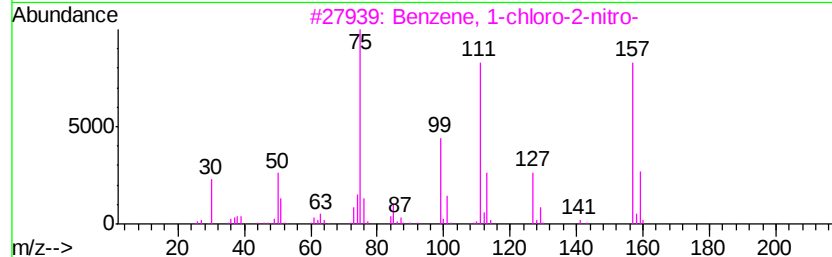
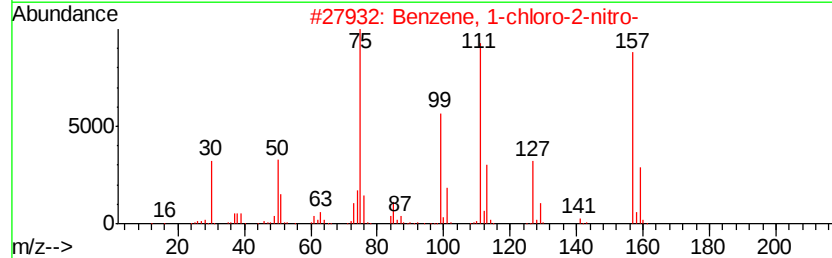
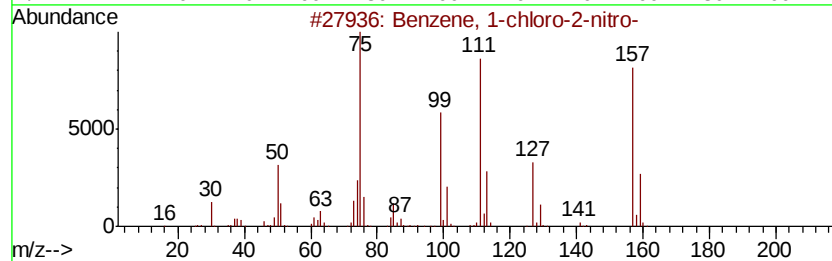
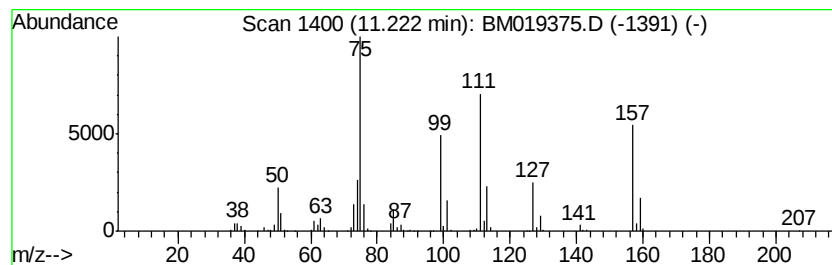
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	30.71 ng/ul	2139390	Naphthalene-d8	10.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
4		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	93



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
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Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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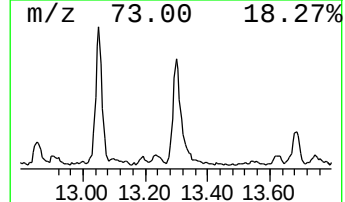
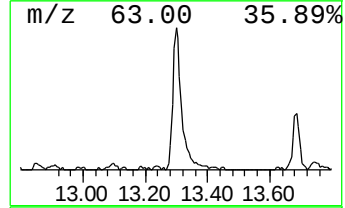
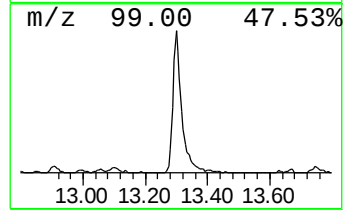
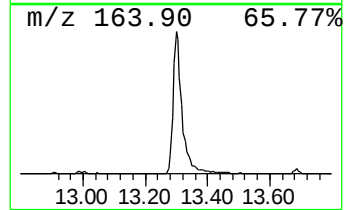
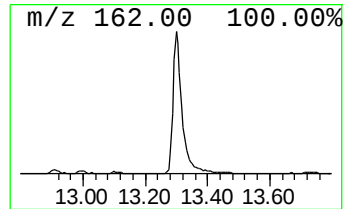
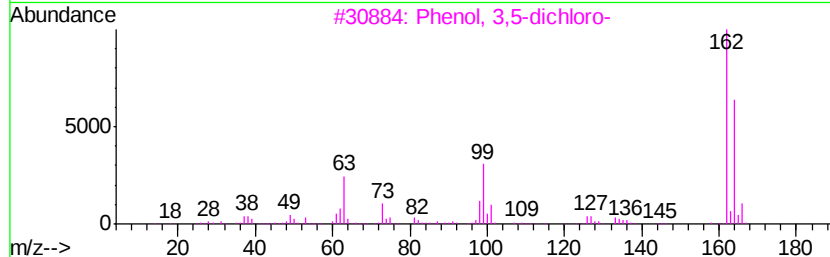
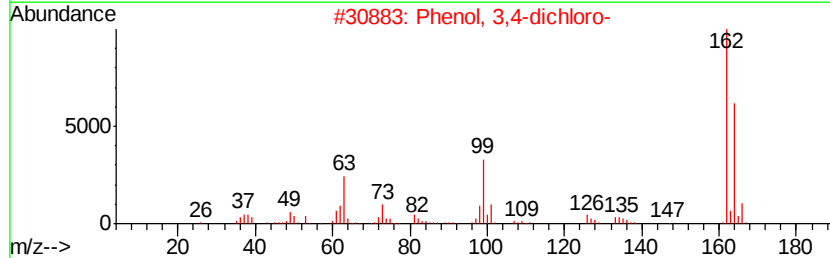
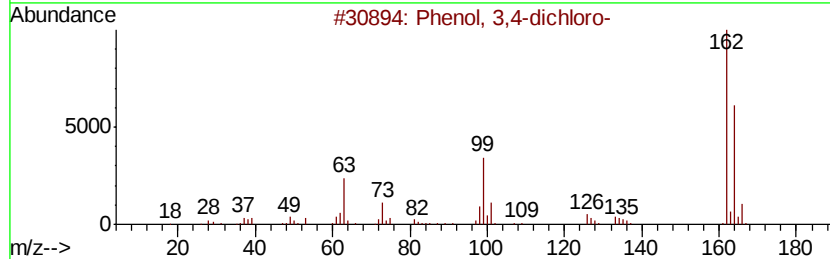
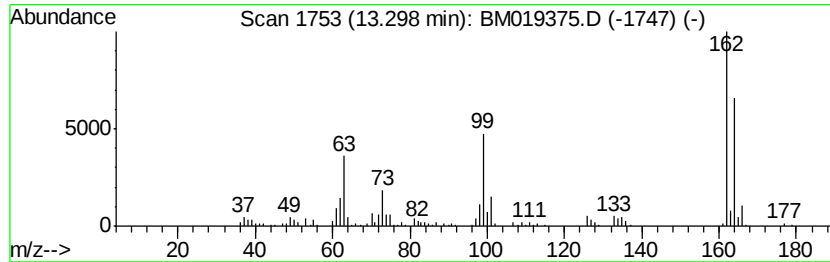
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenol, 3,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.30	6.61 ng/ul	604654	Acenaphthene-d10	14.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dichloro-			162	C6H4Cl2O	000095-77-2	94
2	Phenol, 3,4-dichloro-			162	C6H4Cl2O	000095-77-2	93
3	Phenol, 3,5-dichloro-			162	C6H4Cl2O	000591-35-5	92
4	Phenol, 3,5-dichloro-			162	C6H4Cl2O	000591-35-5	91
5	Phenol, 3,5-dichloro-			162	C6H4Cl2O	000591-35-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM032319\
Data File : BM019375.D
Acq On : 24 Mar 2019 05:00
Operator : JU/SJ
Sample : K2025-14
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM032219MA.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
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Benzene, 1,4-dich...	7.50	4.9	ng/ul	24027600	1	7.62	97190200	20.0
Benzene, 1,3-dich...	7.67	20.0	ng/ul	97190200	1	7.62	97190200	20.0
Benzene, 1,2-dich...	7.99	23.8	ng/ul	115491000	1	7.62	97190200	20.0
Benzene, 1-bromo-...	9.10	3.2	ng/ul	221085	2	10.39	1393200	20.0
Phenol, 2,6-dichl...	10.18	6.5	ng/ul	456314	2	10.39	1393200	20.0
Benzene, 1,3,5-tr...	10.26	404.2	ng/ul	28154500	2	10.39	1393200	20.0
Benzene, 1,2,3-tr...	10.78	175.7	ng/ul	12237700	2	10.39	1393200	20.0
Benzene, 1-chloro...	11.01	115.6	ng/ul	8052450	2	10.39	1393200	20.0
Benzene, 1-chloro...	11.22	30.7	ng/ul	2139390	2	10.39	1393200	20.0
Phenol, 3,4-dichl...	13.30	6.6	ng/ul	604654	3	14.26	1829490	20.0