

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032717\
 Data File : BG026446.D
 Acq On : 27 Mar 2017 23:50
 Operator : SJ/MA
 Sample : I2358-16
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COBMO

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_G\METHODS\SOM-EPA-BG032717MA.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	5.375	376	389	403	rBV	17574091	41181950	44.95%	13.407%
2	7.373	721	729	738	rBV	543021	936378	1.02%	0.305%
3	7.579	756	764	780	rBV2	526487	1250089	1.36%	0.407%
4	7.943	816	826	838	rBV	8474800	17416987	19.01%	5.670%
5	8.119	838	856	861	rBV	21556872	68130101	74.37%	22.180%
6	8.460	896	914	916	rBV4	22265815	91614694	100.00%	29.825%
7	9.206	1033	1041	1043	rBV	565352	979811	1.07%	0.319%
8	9.253	1043	1049	1059	rVB	3799529	7050907	7.70%	2.295%
9	9.923	1158	1163	1182	rVB	492895	924746	1.01%	0.301%
10	10.469	1248	1256	1263	rBV	769942	1519622	1.66%	0.495%
11	10.716	1289	1298	1313	rVV	11796238	23104464	25.22%	7.522%
12	10.845	1313	1320	1328	rVV	816803	1438007	1.57%	0.468%
13	11.233	1377	1386	1398	rBV	5215967	9405272	10.27%	3.062%
14	11.439	1408	1421	1437	rVB	1935170	3394396	3.71%	1.105%
15	11.650	1449	1457	1467	rBV	506889	918061	1.00%	0.299%
16	13.460	1758	1765	1776	rBV	1214196	1987863	2.17%	0.647%
17	14.047	1859	1865	1872	rBV	1581597	2338127	2.55%	0.761%
18	14.341	1908	1915	1922	rBV	1822710	2933378	3.20%	0.955%
19	14.647	1958	1967	1978	rBV2	1229119	2052562	2.24%	0.668%
20	15.634	2128	2135	2144	rBV	2562767	3847546	4.20%	1.253%
21	15.751	2148	2155	2168	rVV	794612	1162140	1.27%	0.378%
22	17.379	2426	2432	2440	rBV2	1695018	2489305	2.72%	0.810%
23	17.479	2442	2449	2459	rBV	2760026	4235624	4.62%	1.379%
24	19.758	2829	2837	2844	rBV	3484735	5234636	5.71%	1.704%
25	21.627	3148	3155	3167	rBV	1898120	3176528	3.47%	1.034%
26	24.588	3648	3659	3677	rVB2	1859427	5188146	5.66%	1.689%
27	24.805	3685	3696	3709	rVB3	1170613	3260365	3.56%	1.061%

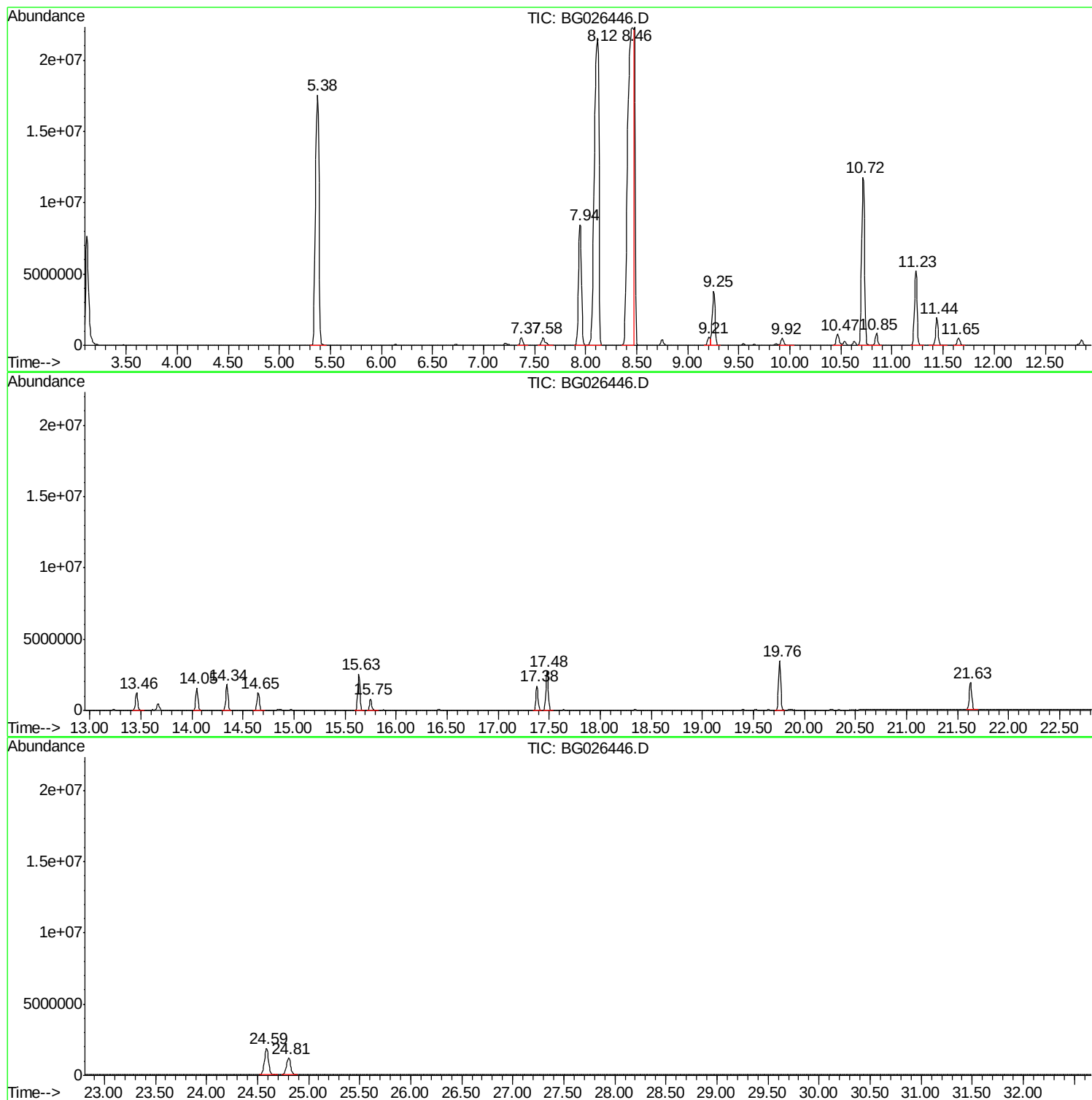
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Instrument :
BNA_G
ClientSampleId :
COBM0

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG032717MA.M
Quant Title : SVOA CALIBRATION

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



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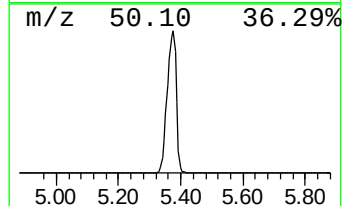
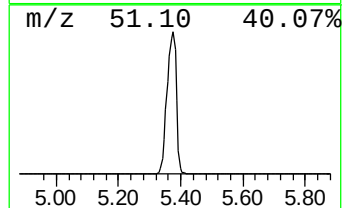
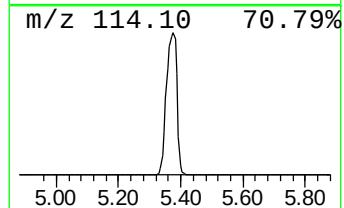
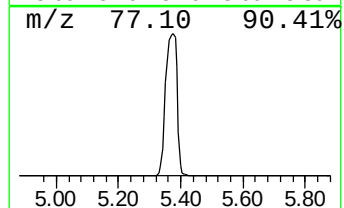
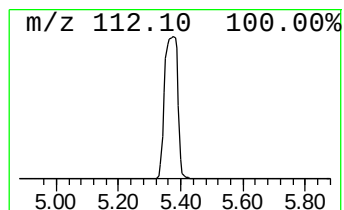
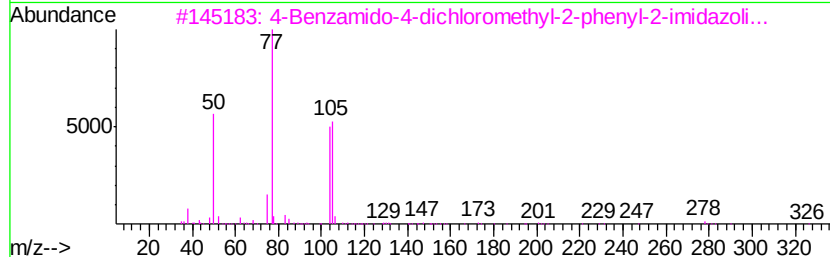
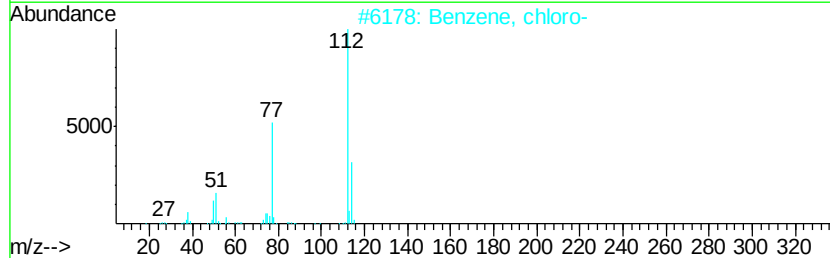
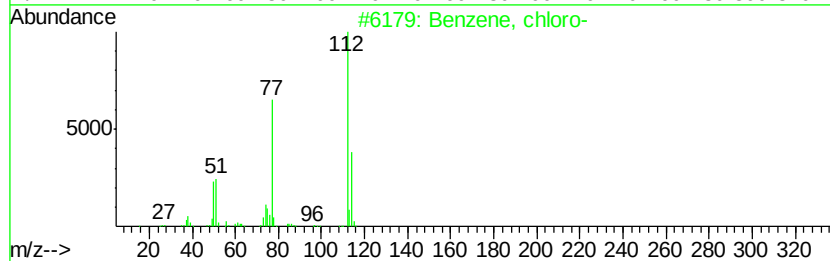
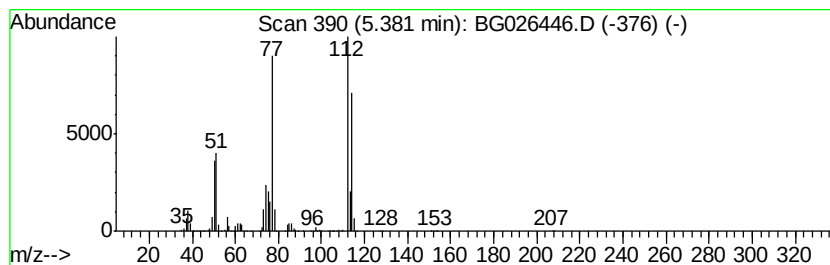
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG032717MA.M
Quant Title : SVOA CALIBRATION

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.
5.38	12.09 ng/ul	41182000	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	87
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	64
3		4-Benzamido-4-dichloromethyl-2-p...	361	C17H13Cl2N3O2	076543-29-8	25
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	12
5		Benzene, chloro-	112	C6H5Cl	000108-90-7	9



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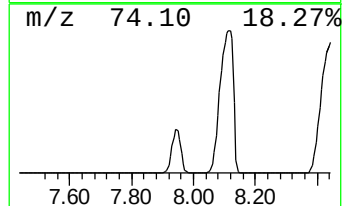
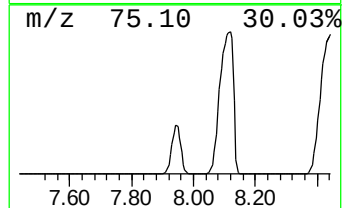
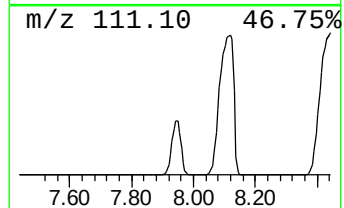
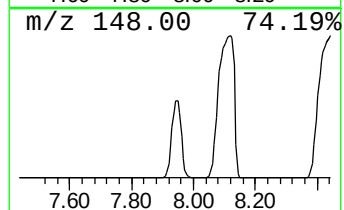
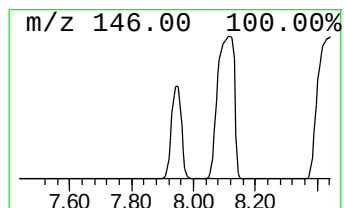
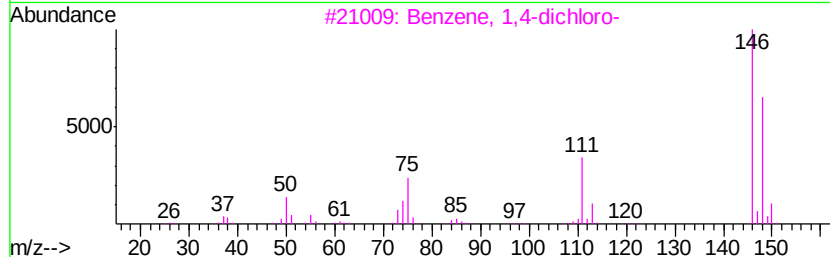
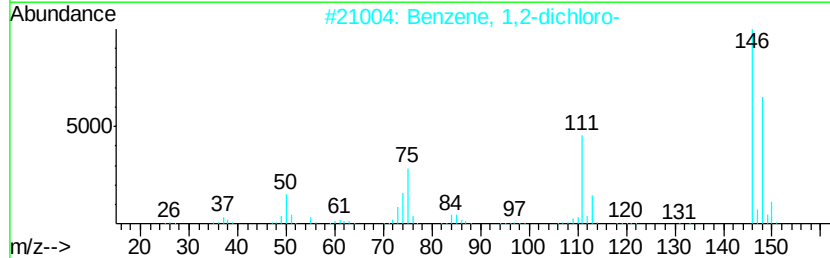
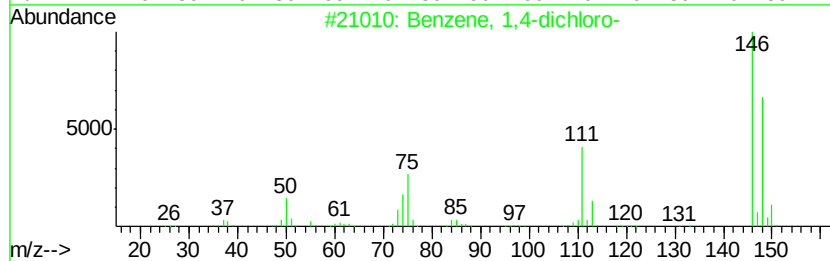
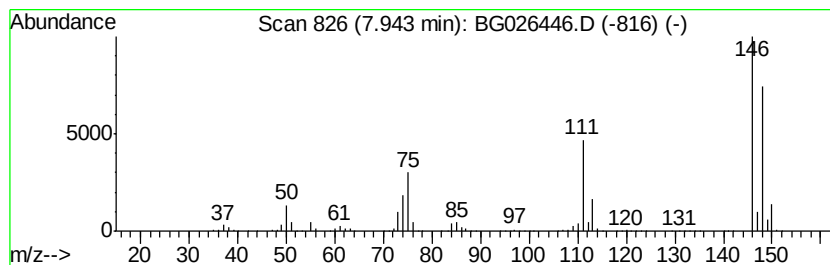
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,4-dichloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.94	5.11 ng/ul	17417000	1,4-Dichlorobenzene-d4	8.06

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	96
3		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95



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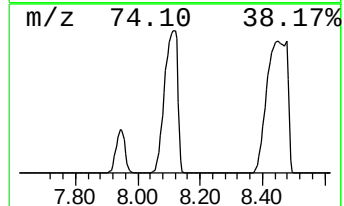
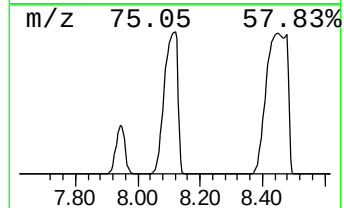
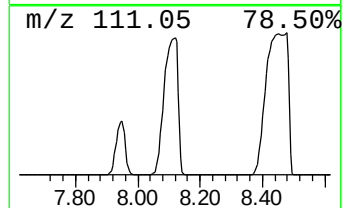
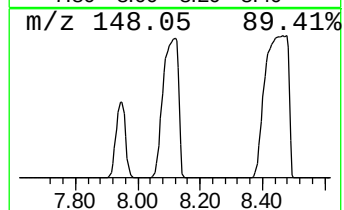
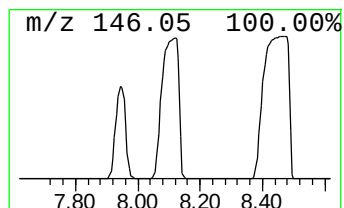
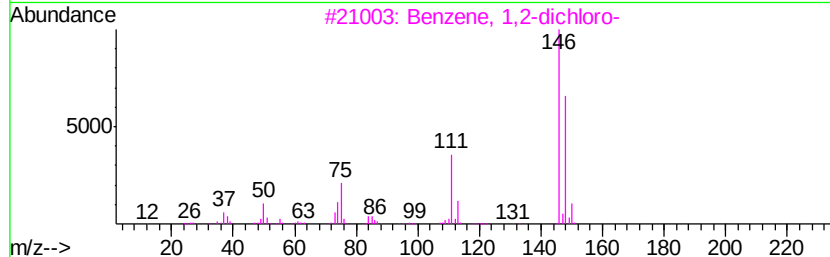
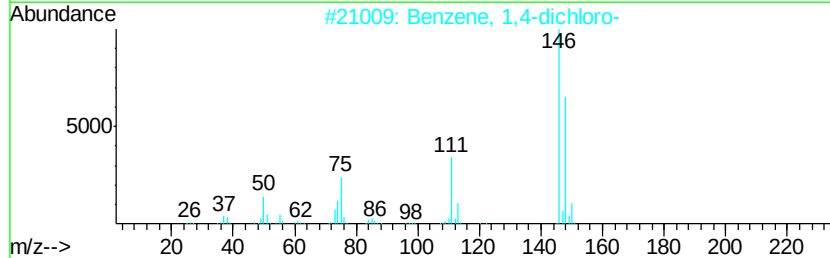
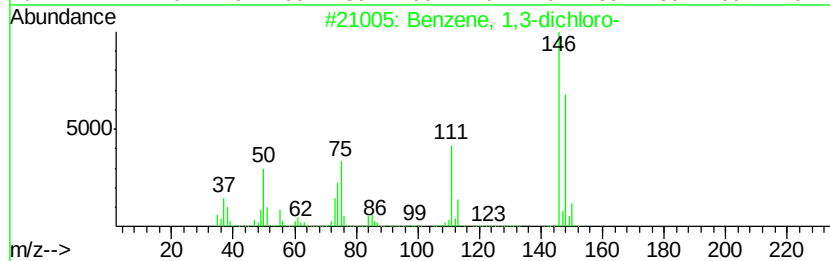
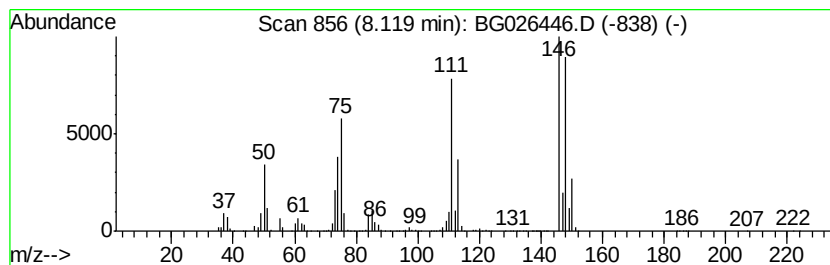
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	20.00 ng/ul	68130100	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	90
3		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	87
4		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	87
5		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	83



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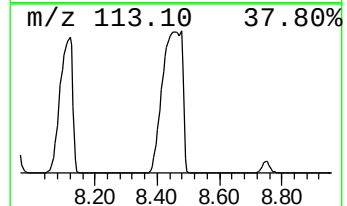
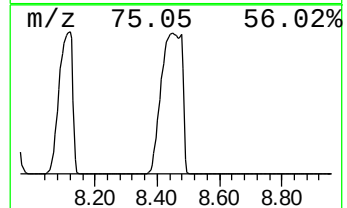
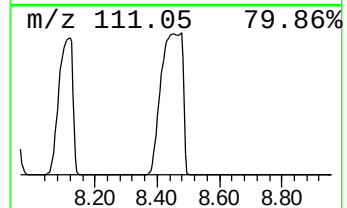
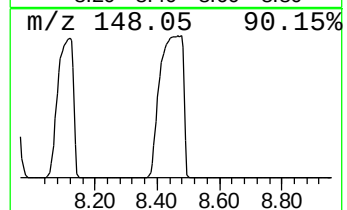
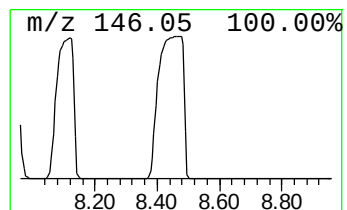
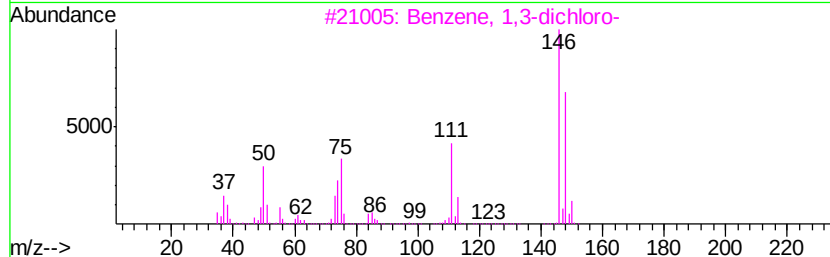
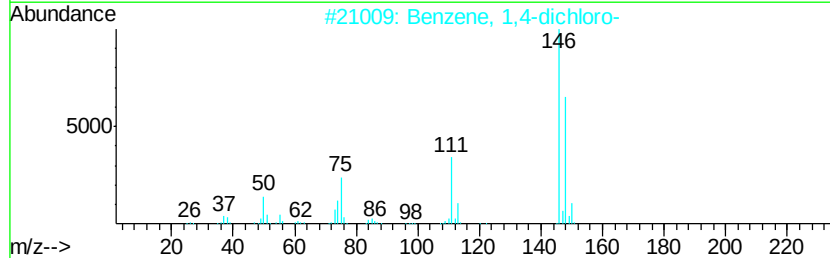
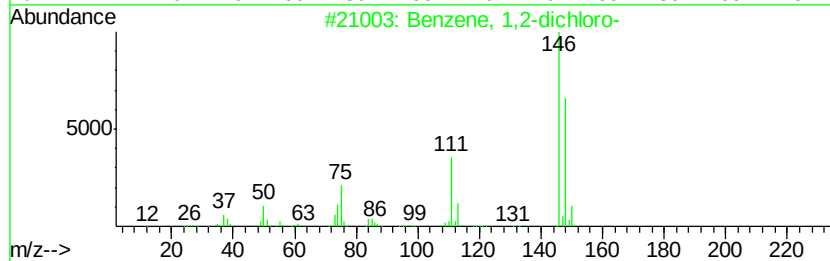
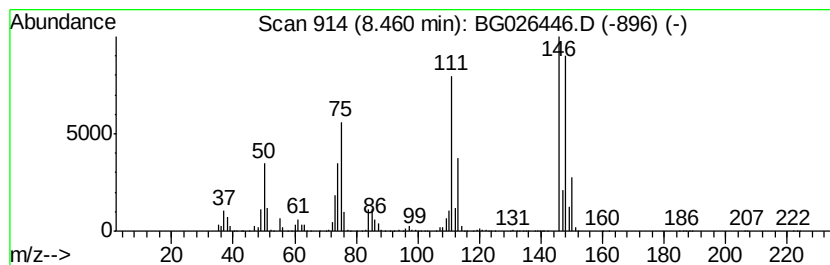
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.46	26.89 ng/ul	91614700	1,4-Dichlorobenzene-d4	8.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	93
2		Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	90
3		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	90
4		Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	81
5		Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	81



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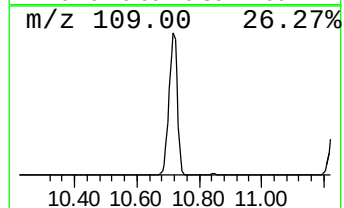
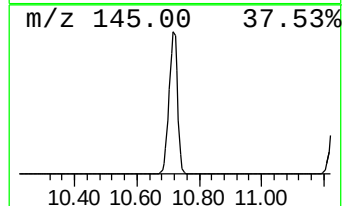
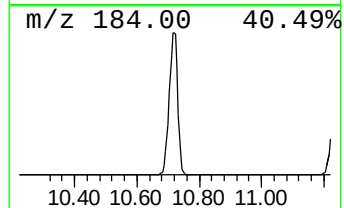
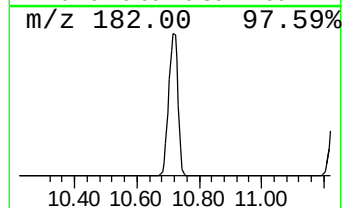
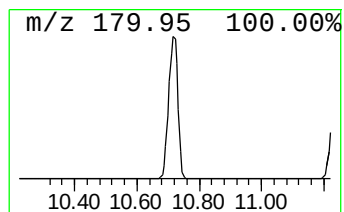
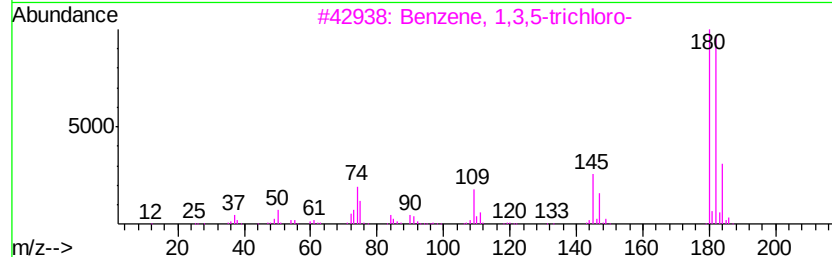
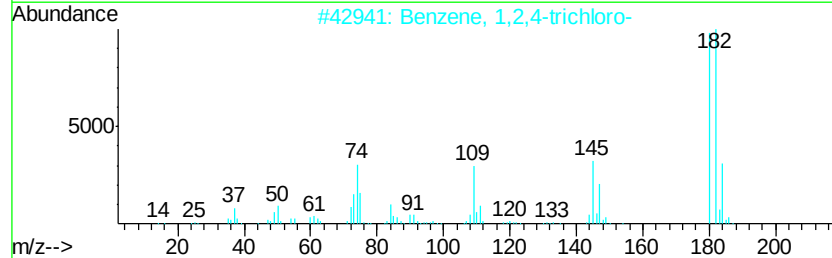
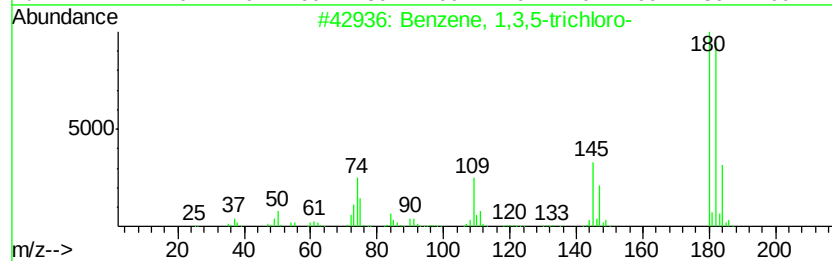
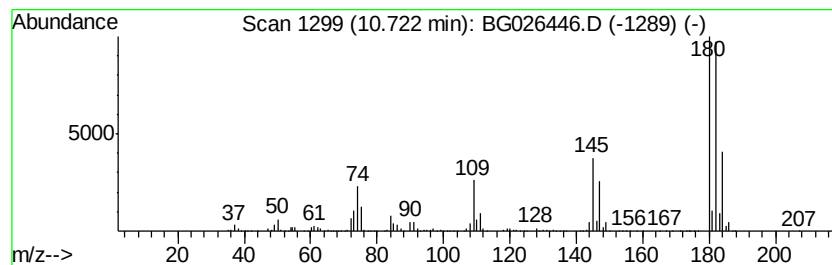
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.72	321.34 ng/ul	23104500	Napthalene-d8	10.85
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3	97
2	Benzene, 1,2,4-trichloro-	180 C6H3Cl3	000120-82-1	97
3	Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3	96
4	Benzene, 1,3,5-trichloro-	180 C6H3Cl3	000108-70-3	96
5	Benzene, 1,2,3-trichloro-	180 C6H3Cl3	000087-61-6	95



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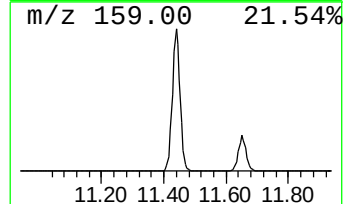
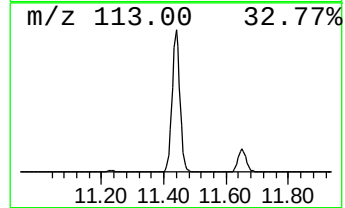
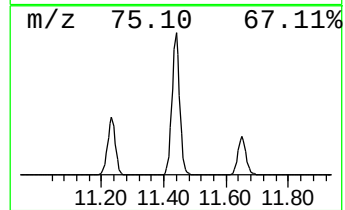
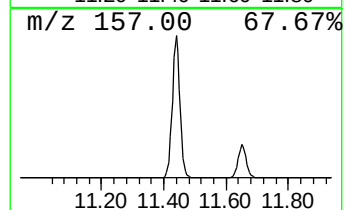
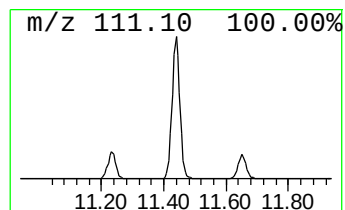
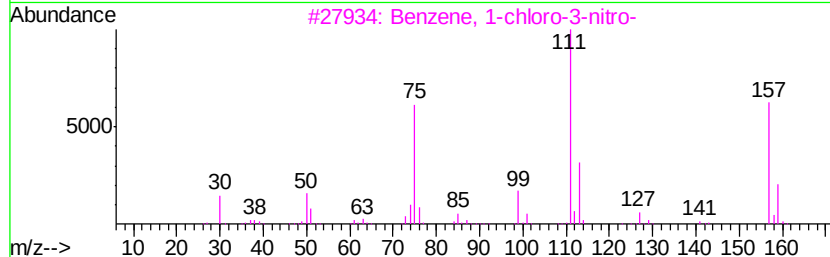
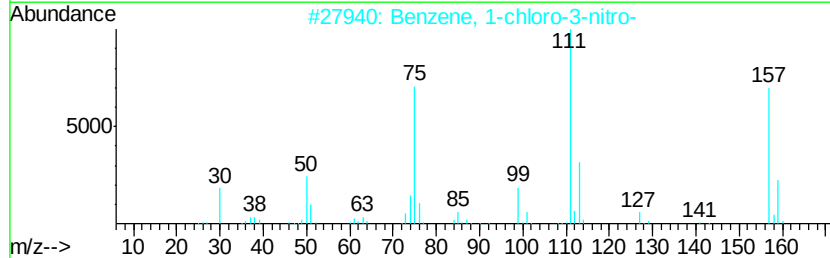
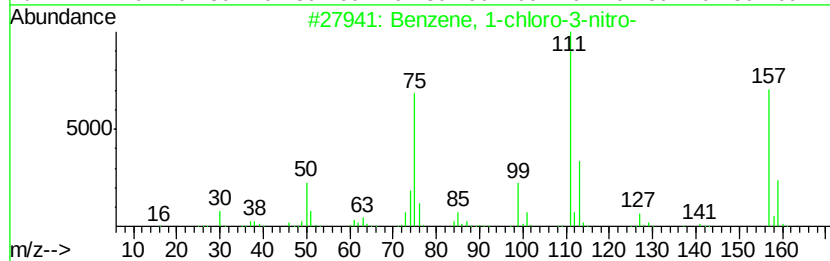
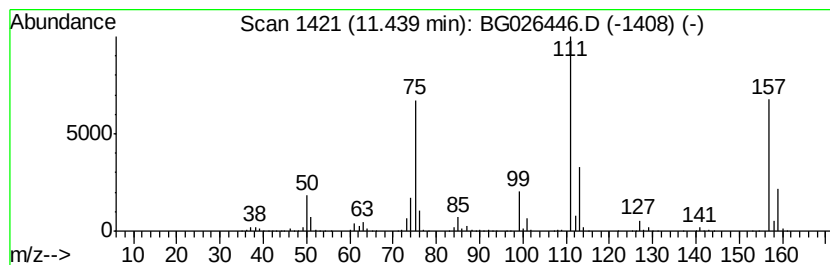
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.44	47.21 ng/ul	3394400	Napthalene-d8	10.85

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-2-nitro-	157	C6H4ClNO2	000088-73-3	87



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032717\
Data File : BG026446.D
Acq On : 27 Mar 2017 23:50
Operator : SJ/MA
Sample : I2358-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

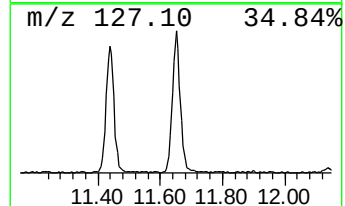
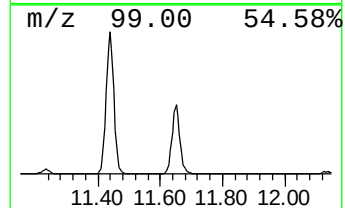
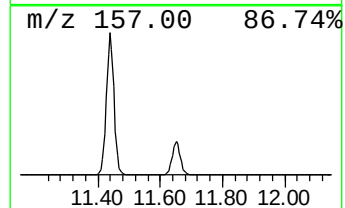
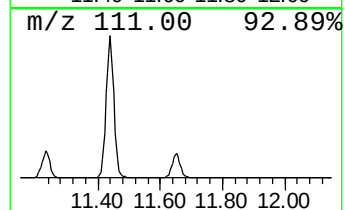
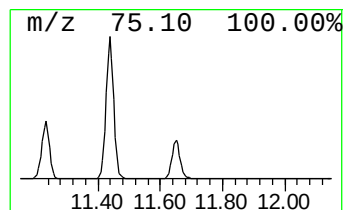
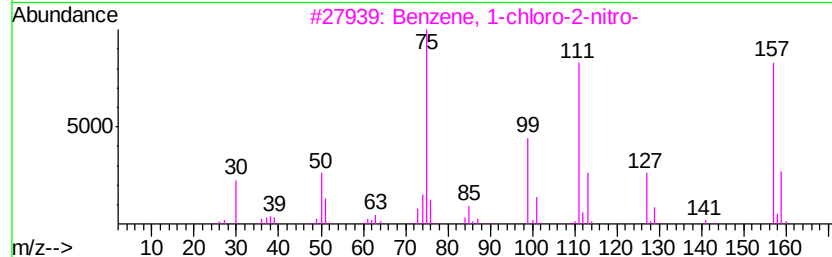
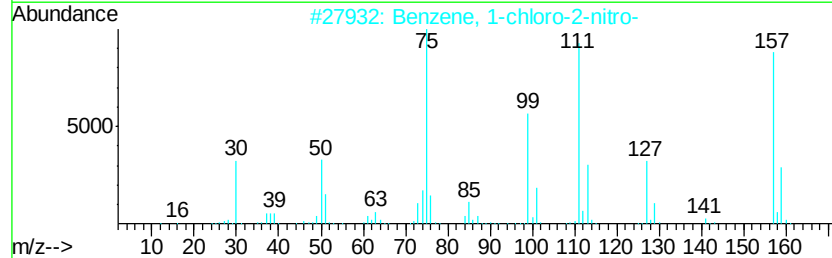
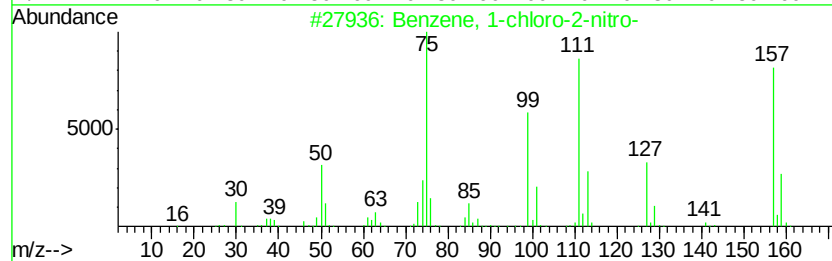
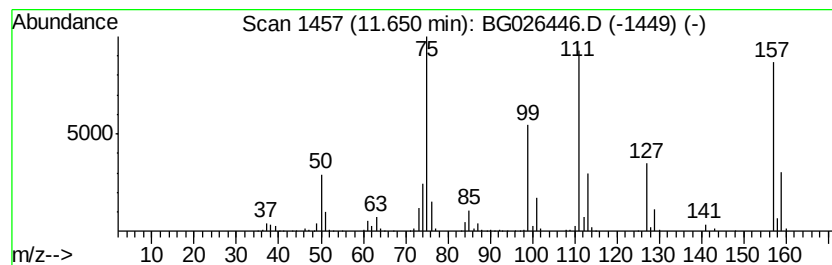
Instrument :
BNA_G
ClientSampleId :
COBMO

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG032717MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1-chloro-2-nitro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.65	12.77 ng/ul	918061	Napthalene-d8	10.85
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	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 98
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5 98
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5 96



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032717\
Data File : BG026446.D
Acq On : 27 Mar 2017 23:50
Operator : SJ/MA
Sample : I2358-16
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0BM0

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG032717MA.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-	5.38	12.1	ng/ul	41182000	1	8.06	68130100	20.0
Benzene, 1,4-dich...	7.94	5.1	ng/ul	17417000	1	8.06	68130100	20.0
Benzene, 1,3-dich...	8.12	20.0	ng/ul	68130100	1	8.06	68130100	20.0
Benzene, 1,2-dich...	8.46	26.9	ng/ul	91614700	1	8.06	68130100	20.0
Benzene, 1,3,5-tr...	10.72	321.3	ng/ul	23104500	2	10.85	1438010	20.0
Benzene, 1-chloro...	11.44	47.2	ng/ul	3394400	2	10.85	1438010	20.0
Benzene, 1-chloro...	11.65	12.8	ng/ul	918061	2	10.85	1438010	20.0