LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0R35

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BM044522.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	30	34	46	rVB	72484	114703	0.10%	0.020%
2	3.687	99	102	120	rBV	347721	598072	0.54%	0.107%
3	5.093	333	341	347	rBV	26324119	82475095	74.65%	14.713%
4	6.428	562	568	574	rBV	305384	473532	0.43%	0.084%
5	6.716	613	617	623	rVB	94949	158385	0.14%	0.028%
6	6.910	644	650	661	rBV2	135885	267703	0.24%	0.048%
7	7.093	672	681	692	rBV	1058681	1905757	1.72%	0.340%
8	7.275	705	712	723	rBV	474043	867541	0.79%	0.155%
9	7.463	738	744	755	rVB	73297	153453	0.14%	0.027%
10	7.640	762	774	785	rBV	22278507	59048938	53.44%	10.534%
11	7.787	785	799	807	rBV	23902006	99158837	89.75%	17.690%
12	7.851	809	810	812	rVB	21040557	8288853	7.50%	1.479%
13	8.110	843	854	863	rBV	23728763	110488884	100.00%	19.711%
14	8.440	903	910	923	rVB	435167	729705	0.66%	0.130%
15	8.898	980	988	991	rBV	1241134	2135823	1.93%	0.381%
16	8.940	991	995	1003	rVB	2860103	4556966	4.12%	0.813%
17	9.222	1035	1043	1049	rBV	858802	1411440	1.28%	0.252%
18	9.439	1071	1080	1084	rBV	198112	334041	0.30%	0.060%
19	9.487	1084	1088	1091	rVV	193195	357451	0.32%	0.064%
20	9.528	1091	1095	1102	rVB	430163	762889	0.69%	0.136%
21	9.610	1102	1109	1114	rBV	459139	821880	0.74%	0.147%
22	9.657	1114	1117	1123	rVB	121078	184703	0.17%	0.033%
23	9.892	1150	1157	1164	rVB	70633	125864	0.11%	0.022%
24	10.139	1192	1199	1207	rBV	648244	1129204	1.02%	0.201%
25	10.398	1231	1243	1247	rBV	23101401	73100579	66.16%	13.041%
26	10.528	1258	1265	1271	rBV	1206464	1879776	1.70%	0.335%
27	10.663	1280	1288	1299	rBV	1083790	1855649	1.68%	0.331%
28	10.910	1318	1330	1339	rBV	18408250	37136676	33.61%	6.625%
29	11.122	1358	1366	1374	rBV	1672194	2713883	2.46%	0.484%
30	11.333	1393	1402	1412	rBV2	355754	665577	0.60%	0.119%
31	11.810	1471	1483	1492	rBV5	78622	205730	0.19%	0.037%
32	12.045	1515	1523	1537	rBV	315240	638491	0.58%	0.114%
33	12.510	1594	1602	1603	rBV2	373868	513623	0.46%	0.092%
34	12.545	1603	1608	1615	rVV	2607287	4140587	3.75%	0.739%
35	12.792	1644	1650	1657	rBV	77286	145542	0.13%	0.026%
36	13.157	1699	1712	1723	rBV	6775474	10415722	9.43%	1.858%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Title : SVOA CALIBRATION

37	13.786	1808	1819	1825	rBV	2625719	3838712	3.47%	0.685%
38	14.063	1858	1866	1875	rBV	3004842	4614663	4.18%	0.823%
39	14.363	1910	1917	1932	rBV	1704810	2561656	2.32%	0.457%
40	14.680	1965	1971	1976	rVB	249405	369456	0.33%	0.066%
41	14.857	1996	2001	2010	rBV2	73312	119601	0.11%	0.021%
42	15.363	2080	2087	2093	rBV	3994636	5691998	5.15%	1.015%
43	15.486	2103	2108	2119	rVV	517756	765859	0.69%	0.137%
44	17.115	2379	2385	2396	rBV	2007248	2890178	2.62%	0.516%
45	17.215	2396	2402	2414	rVV	4378140	6173941	5.59%	1.101%
46	18.021	2532	2539	2546	rBV2	72930	121318	0.11%	0.022%
47	18.492	2614	2619	2627	rVB	872119	1174211	1.06%	0.209%
48	18.986	2696	2703	2709	rVB	213387	291949	0.26%	0.052%
49	19.515	2786	2793	2802	rBV	5194756	7219532	6.53%	1.288%
50	21.039	3048	3052	3060	rBV3	67672	119011	0.11%	0.021%
51	21.303	3092	3097	3105	rBV	2254785	3033841	2.75%	0.541%
52	21.409	3111	3115	3120	rBV	681617	782453	0.71%	0.140%
53	23.415	3448	3456	3464	rBV	3626165	6780586	6.14%	1.210%
54	23.556	3473	3480	3492	rVB	1524712	3011158	2.73%	0.537%
55	23.821	3520	3525	3534	rVB	570816	1019453	0.92%	0.182%

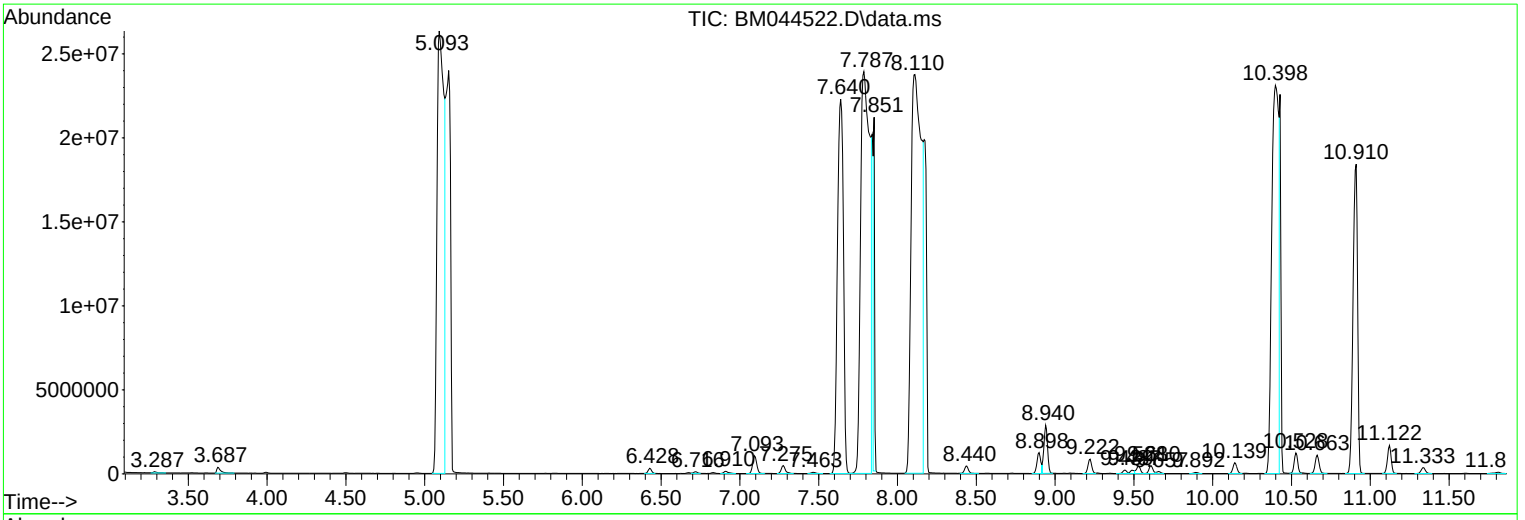
Sum of corrected areas: 560541130

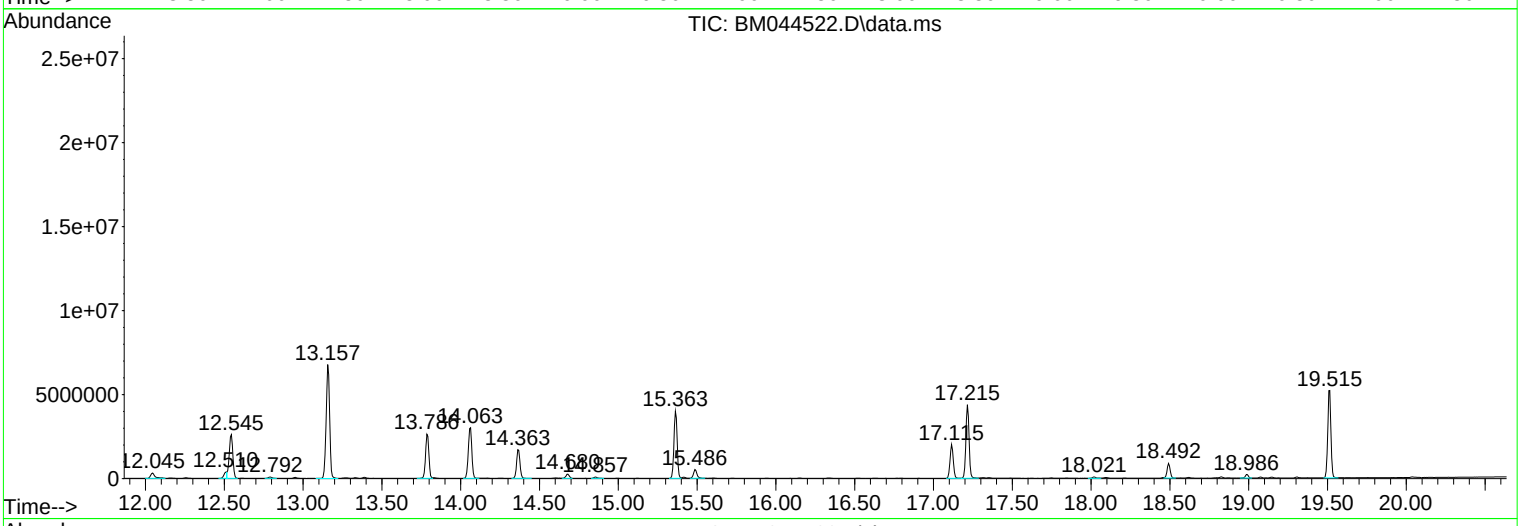
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

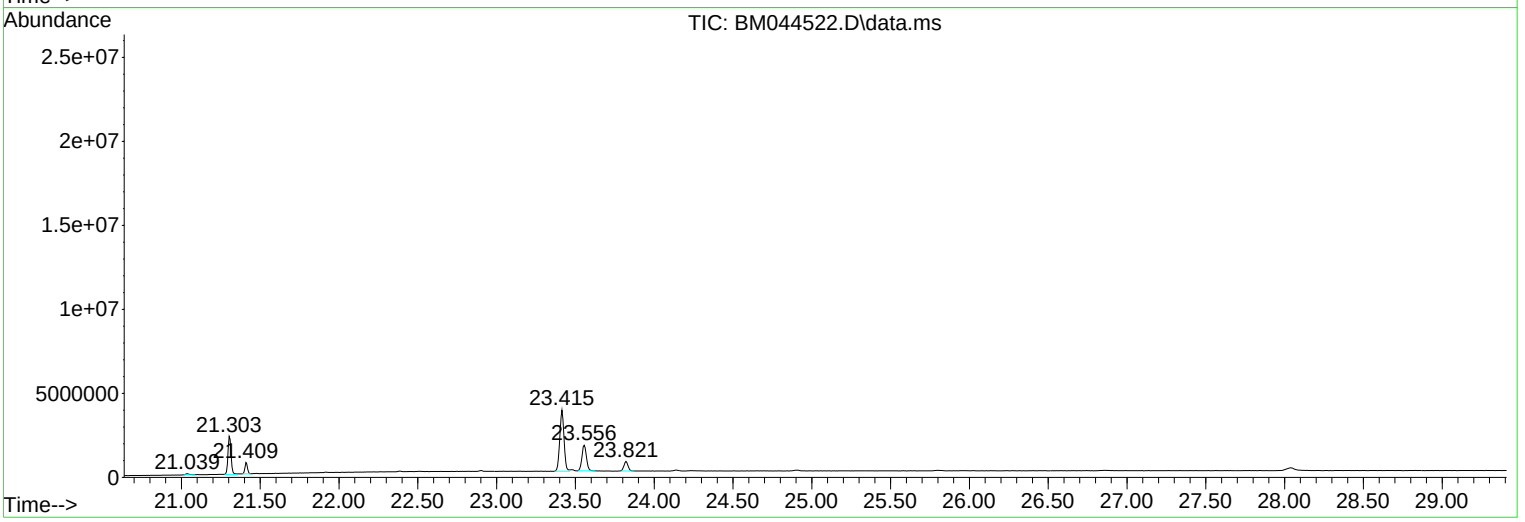
Instrument :
BNA_M
ClientSampleId :
C0R35

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P







Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

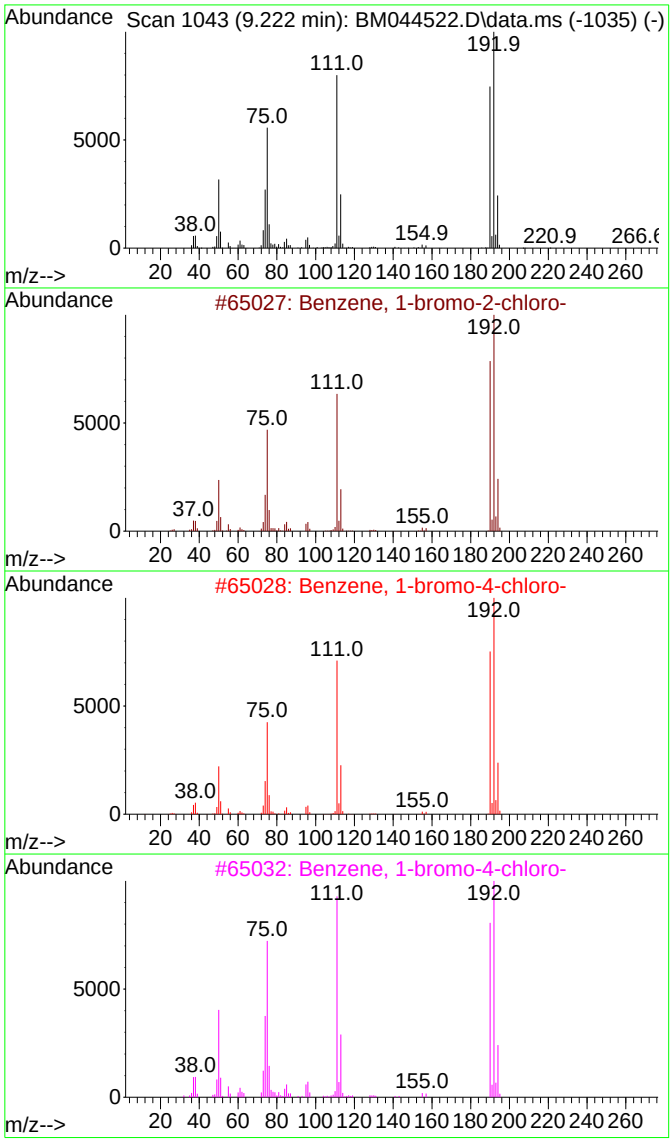
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

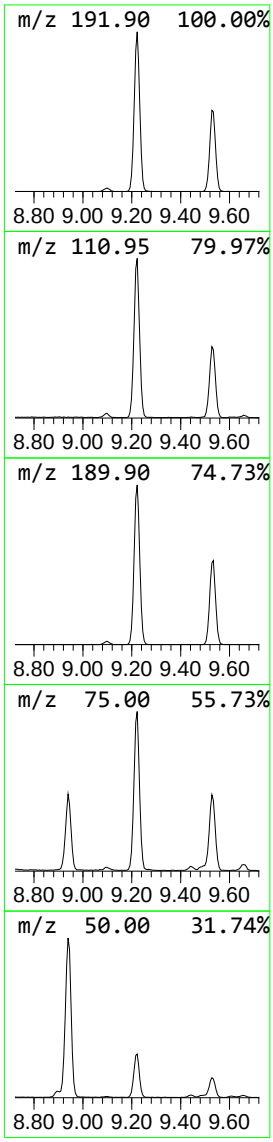
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.222	15.02 ng/ul	1411440	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
4	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

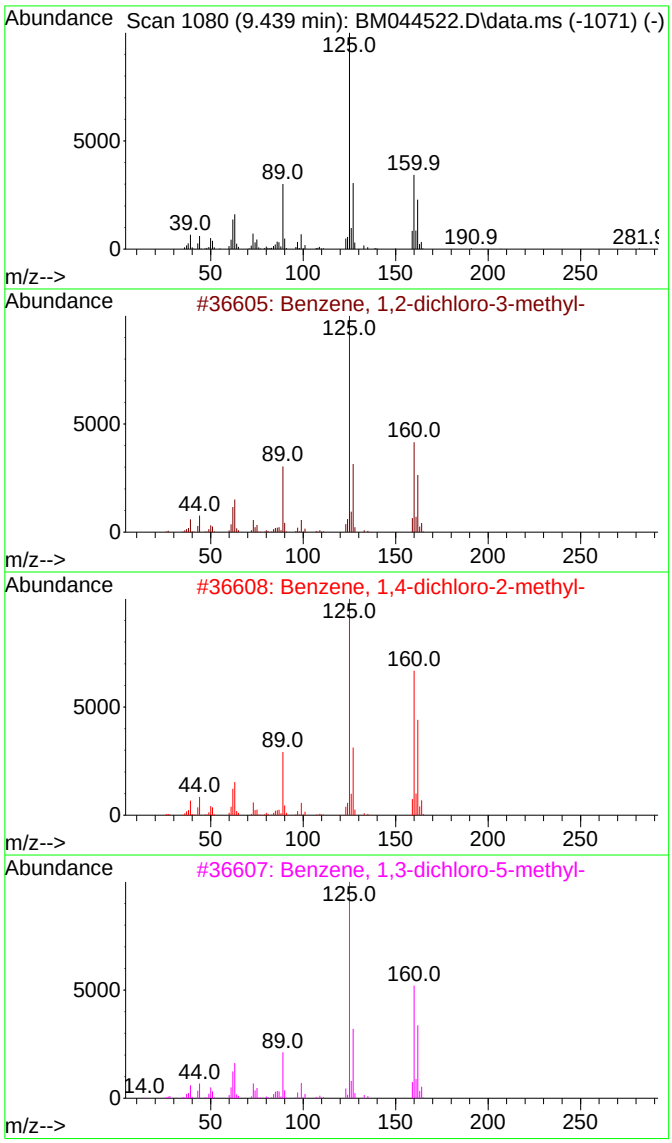
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

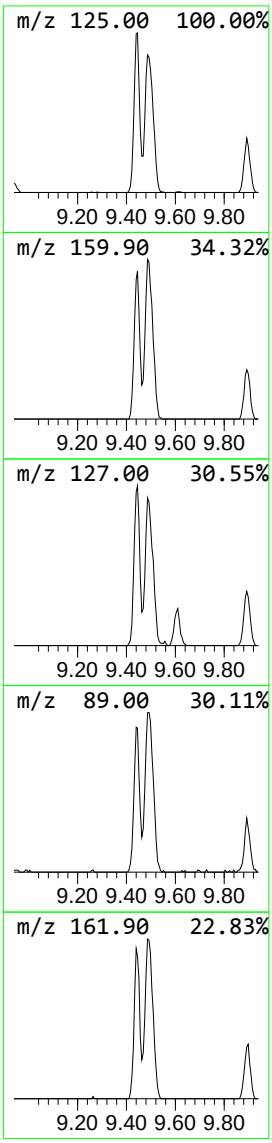
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,2-dichloro-3-met... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.439	3.55 ng/ul	334041	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2	Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	97
3	Benzene, 1,3-dichloro-5-methyl-	160	C7H6Cl2	025186-47-4	97
4	Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
5	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

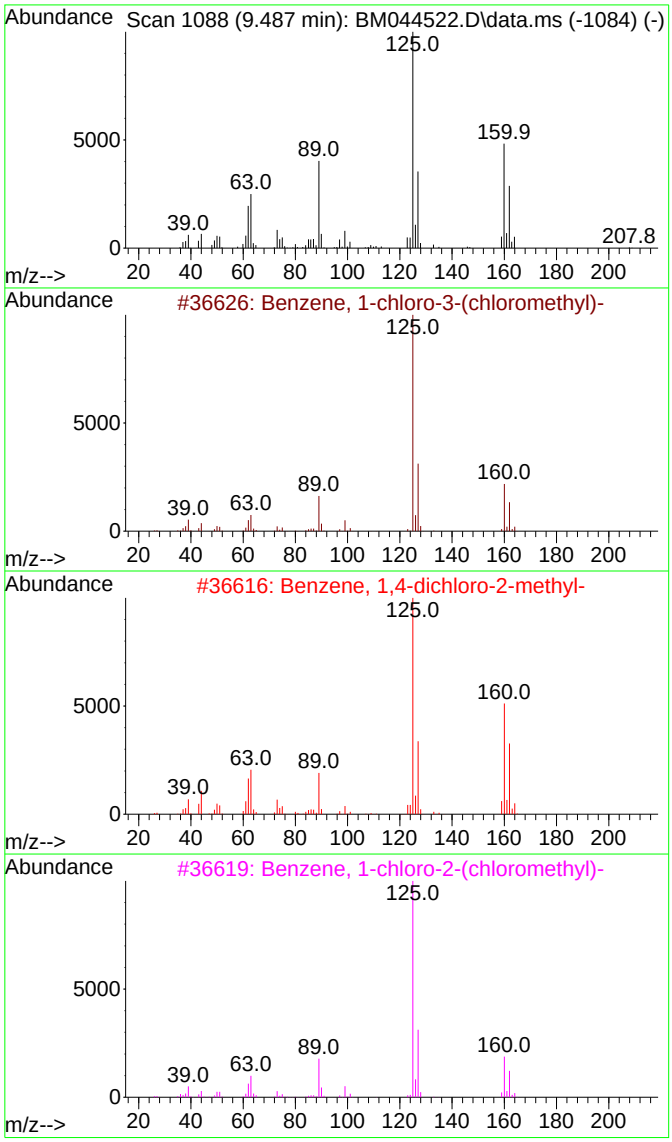
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

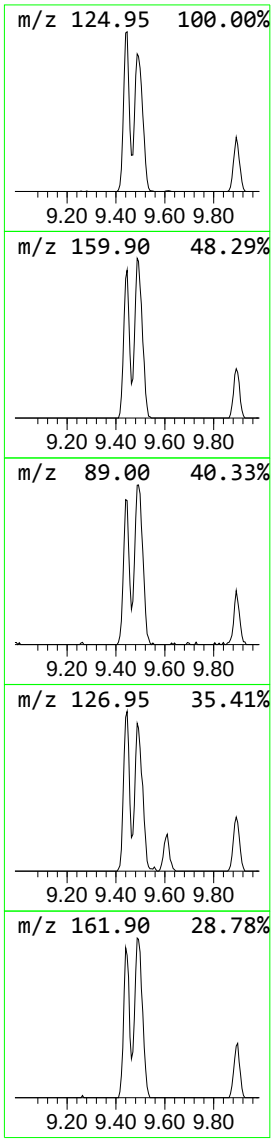
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-chloro-3-(chloro... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.487	3.80 ng/ul	357451	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-(chloromethyl)-	160	C7H6Cl2	000620-20-2	96
2	Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	96
3	Benzene, 1-chloro-2-(chloromethyl)-	160	C7H6Cl2	000611-19-8	95
4	Benzene, 1-chloro-4-(chloromethyl)-	160	C7H6Cl2	000104-83-6	95
5	Benzene, 1,4-dichloro-2-methyl-	160	C7H6Cl2	019398-61-9	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

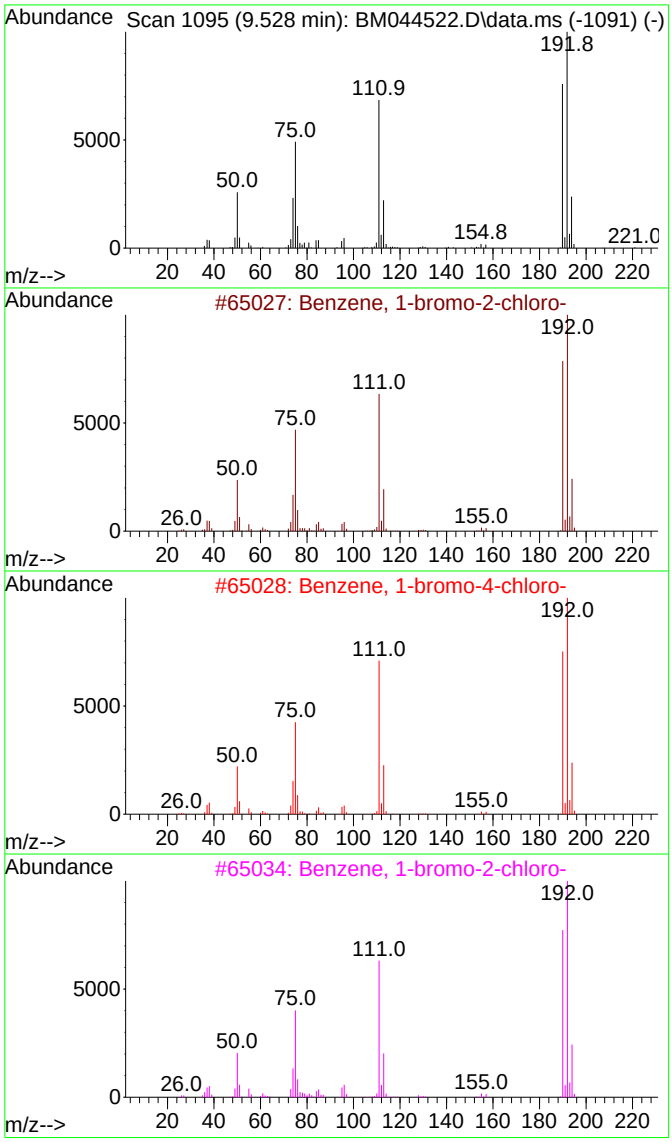
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

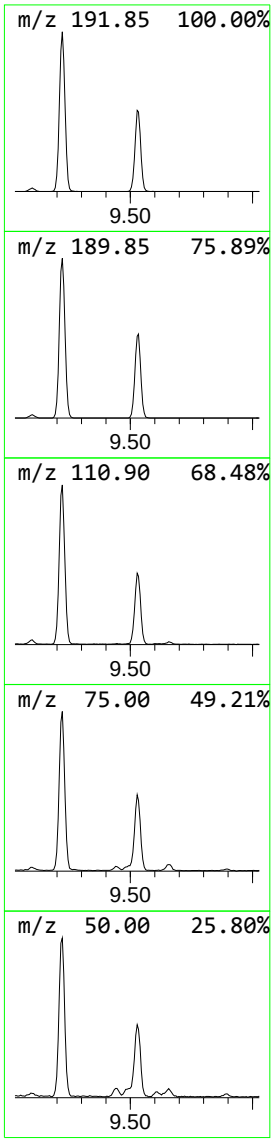
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.528	8.12 ng/ul	762889	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
4	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96
5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

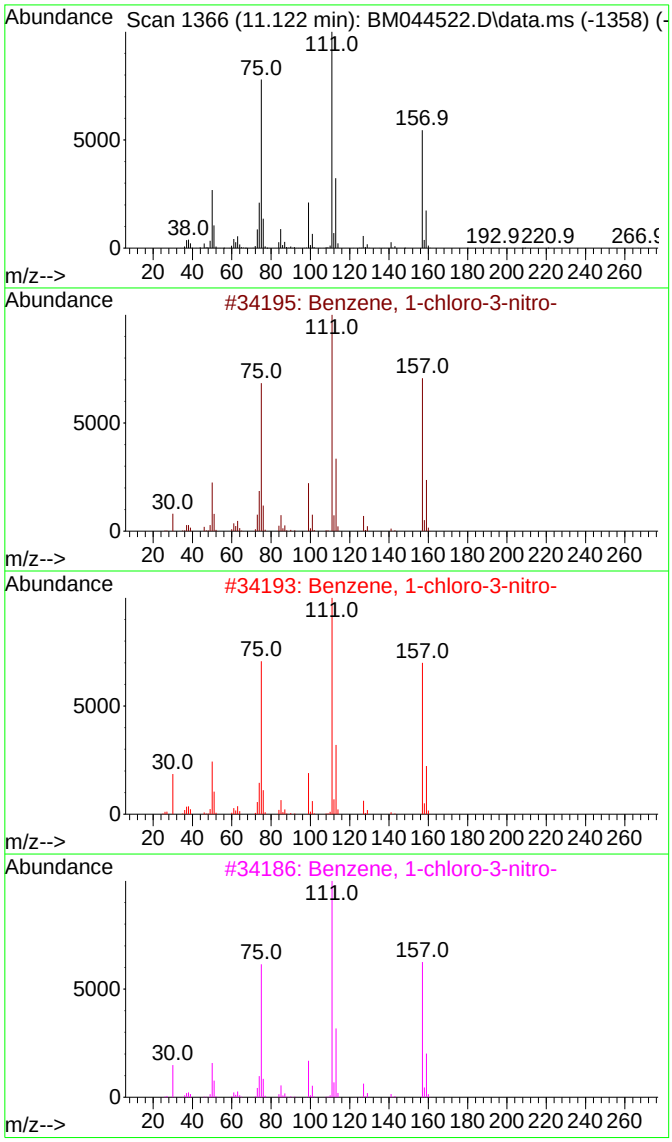
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

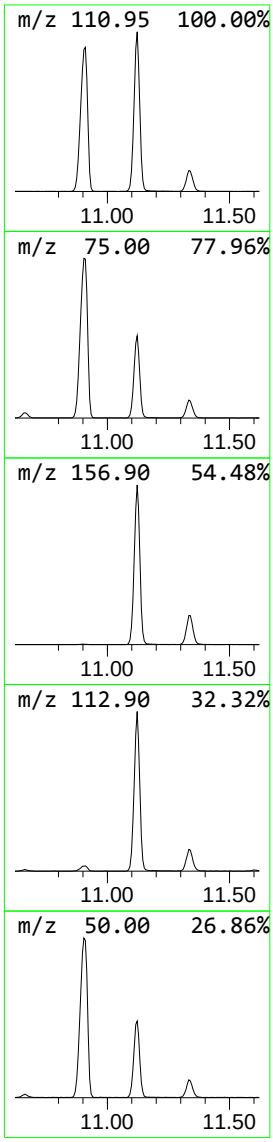
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.122	28.87 ng/ul	2713880	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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	94
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

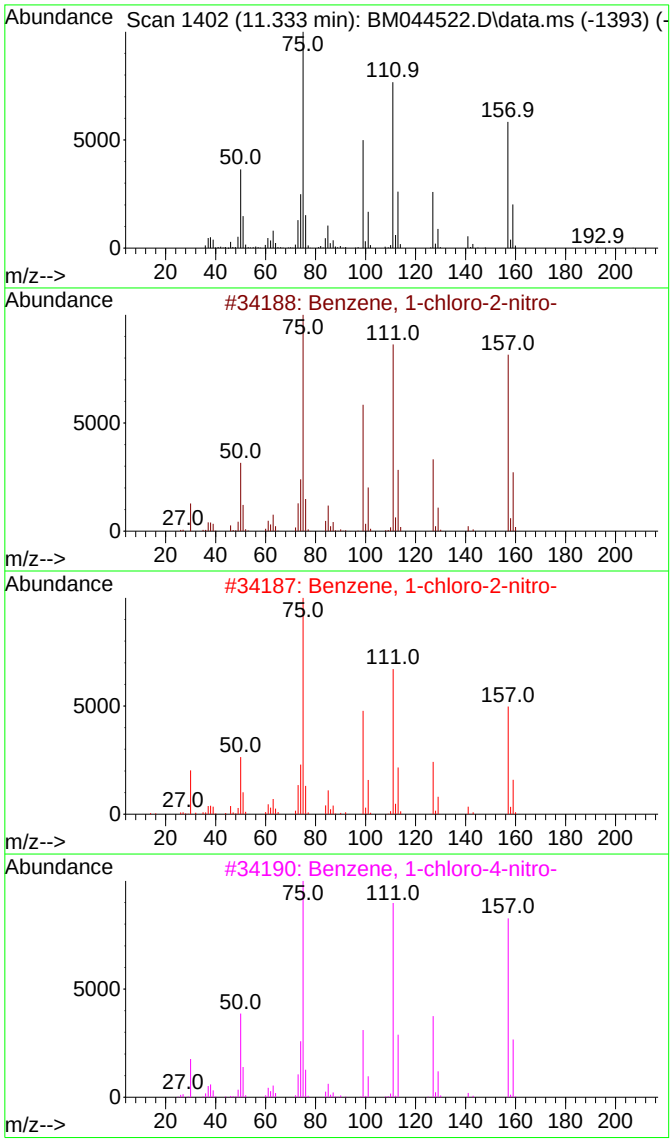
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

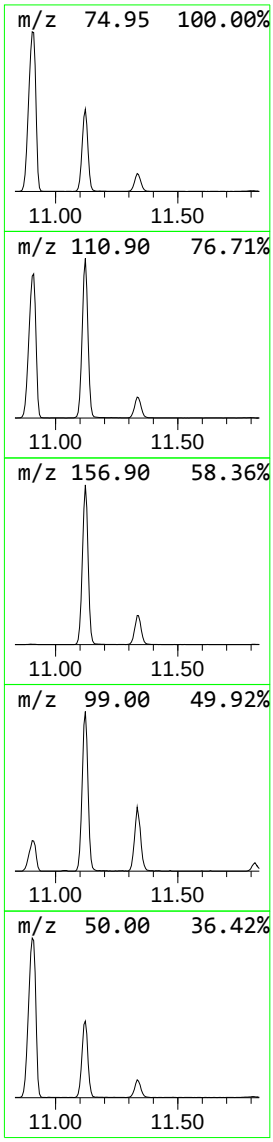
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-chloro-2-nitro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.334	7.08 ng/ul	665577	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

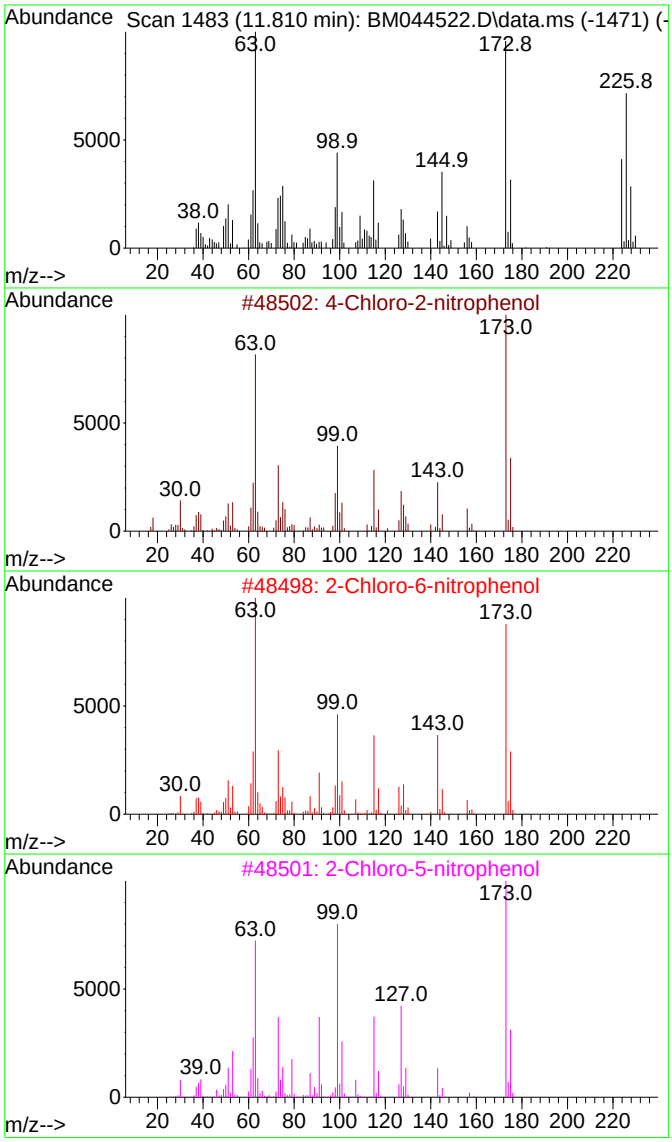
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

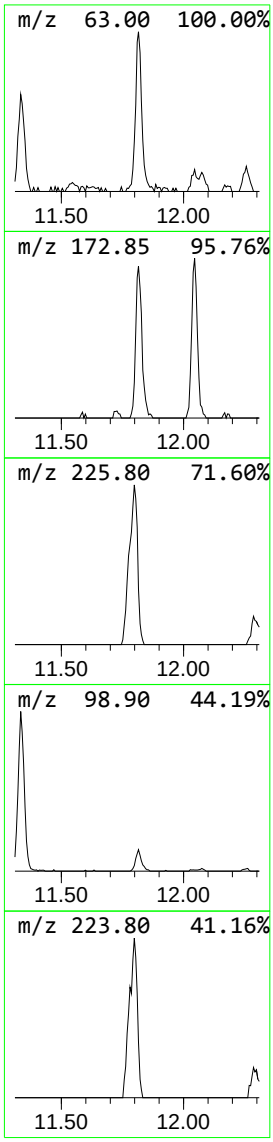
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 4-Chloro-2-nitrophenol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.810	2.19 ng/ul	205730	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	95
2	2-Chloro-6-nitrophenol	173	C6H4ClNO3	000603-86-1	90
3	2-Chloro-5-nitrophenol	173	C6H4ClNO3	000619-10-3	68
4	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	64
5	Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

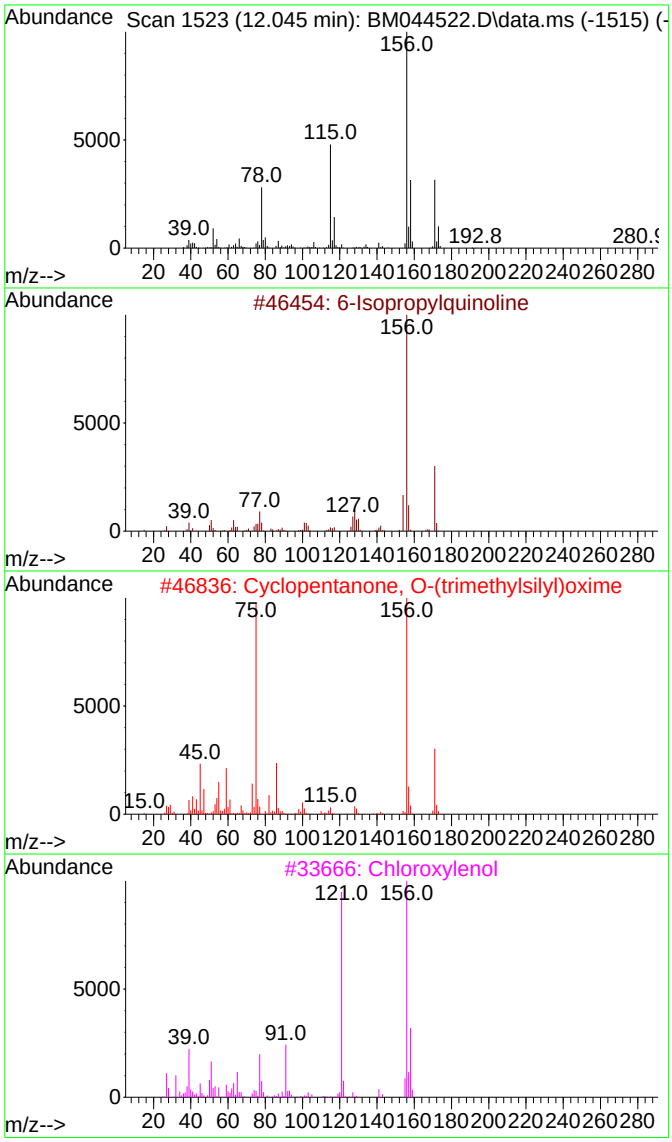
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

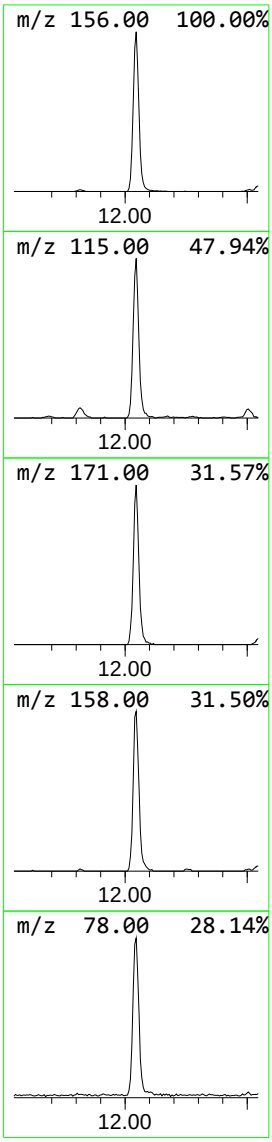
****

Peak Number 13 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.045	6.79 ng/ul	638491	Naphthalene-d8	10.528

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Isopropylquinoline	171	C12H13N	000135-79-5	38
2	Cyclopentanone, O-(trimethylsily...	171	C8H17NOSi	026208-87-7	32
3	Chloroxylenol	156	C8H9ClO	000088-04-0	32
4	Benzene, 1-chloro-4-methoxy-2-me...	156	C8H9ClO	013334-71-9	32
5	Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

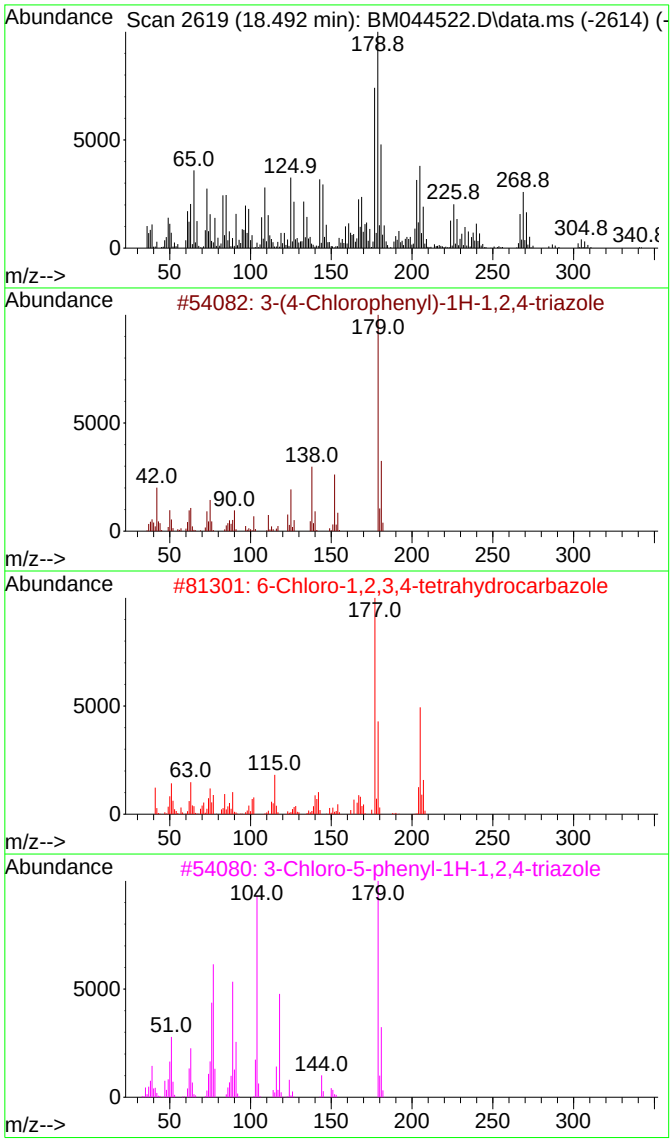
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

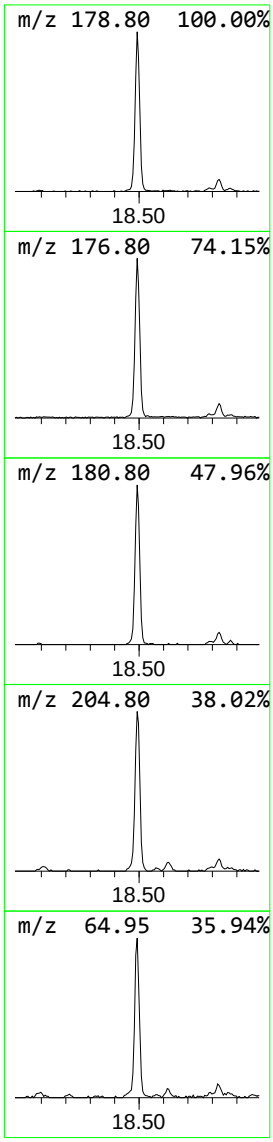
****

Peak Number 14 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.492	8.13 ng/ul	1174210	Phenanthrene-d10	17.115

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-(4-Chlorophenyl)-1H-1,2,4-tria...	179	C8H6ClN3	023195-59-7	35
2	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	35
3	3-Chloro-5-phenyl-1H-1,2,4-triazole	179	C8H6ClN3	031803-05-1	27
4	Benzoic acid, 2-chloro-5-tetrazo...	238	C9H7ClN4O2	1010310-73-5	27
5	Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

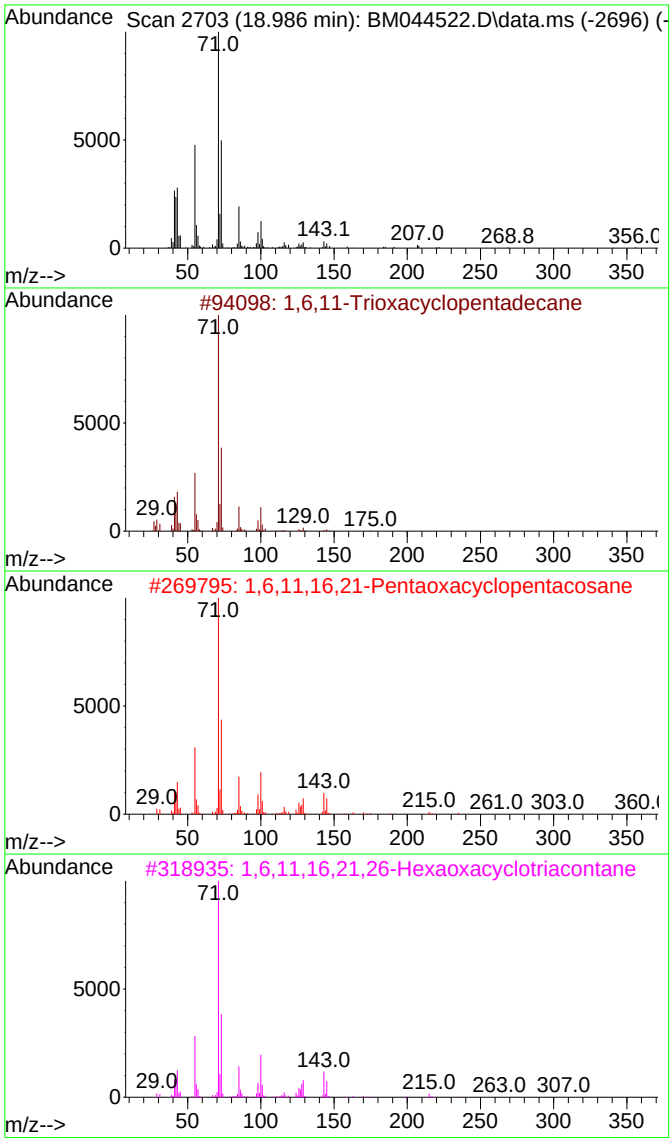
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

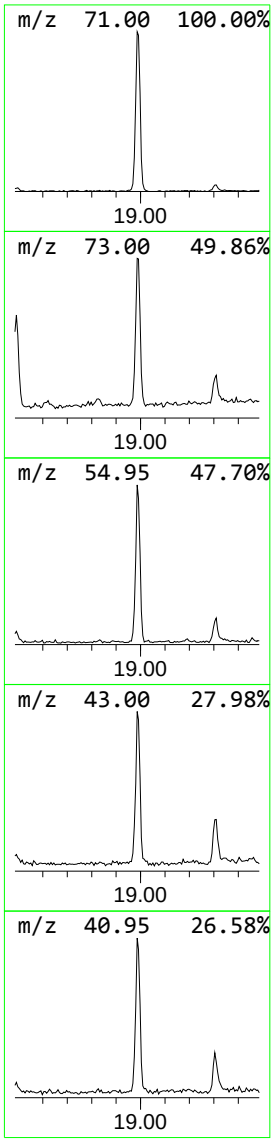
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 1,6,11-Trioxacyclopentadecane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.986	2.02 ng/ul	291949	Phenanthrene-d10	17.115

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,11-Trioxacyclopentadecane	216	C12H24O3	000295-63-6	80
2	1,6,11,16,21-Pentaoxacyclopentac...	360	C20H40O5	056890-57-4	80
3	1,6,11,16,21,26-Hexaoxacyclotria...	432	C24H48O6	064001-05-4	59
4	1-Penten-3-ol, 2-methyl-	100	C6H12O	002088-07-5	53
5	Glutaric acid, ditetrahydrofurfu...	300	C15H24O6	1000359-67-4	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

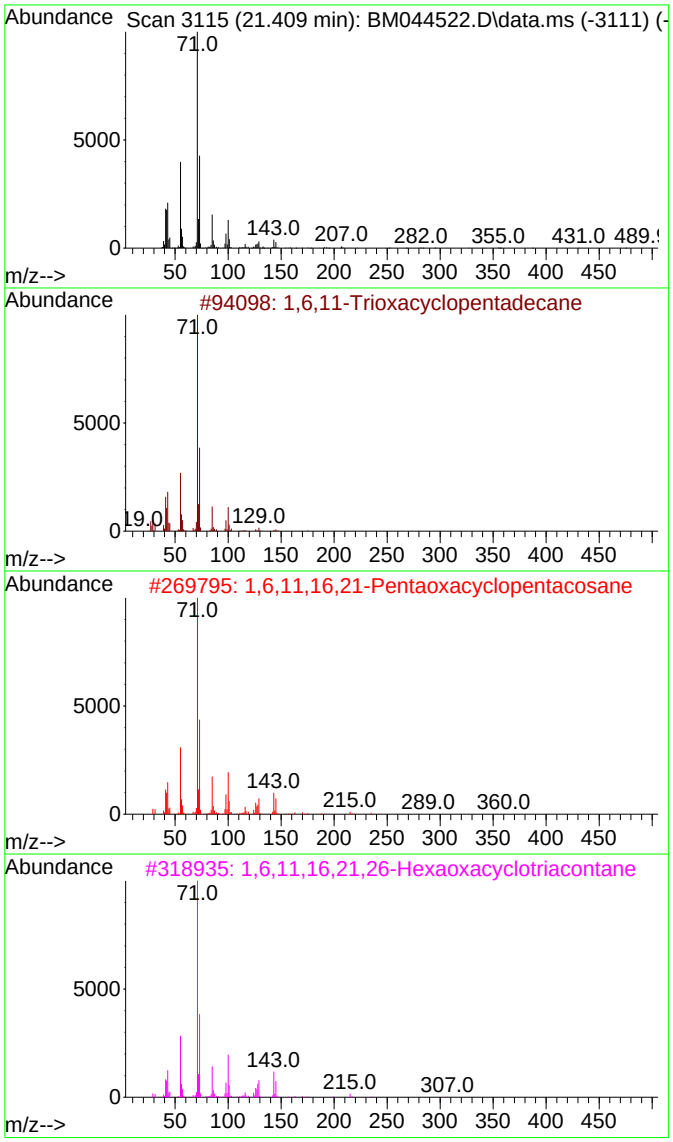
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

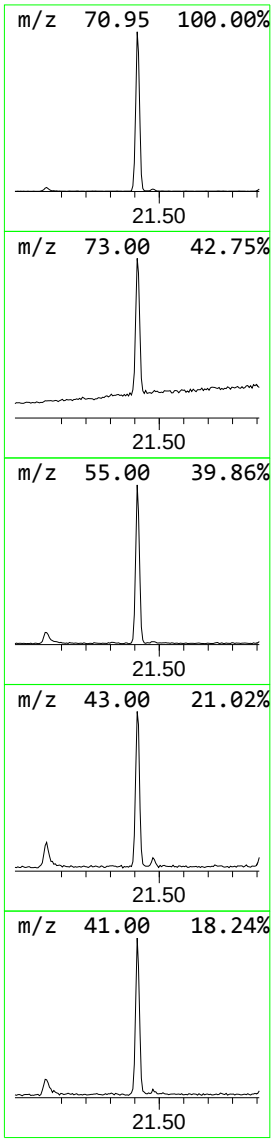
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 1,6,11,16,21-Pentaoxacyclop... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.409	5.16 ng/ul	782453	Chrysene-d12	21.303

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,11-Trioxacyclopentadecane	216	C12H24O3	000295-63-6	91
2	1,6,11,16,21-Pentaoxacyclopentac...	360	C20H40O5	056890-57-4	87
3	1,6,11,16,21,26-Hexaoxacyclotria...	432	C24H48O6	064001-05-4	86
4	1,6,11,16,21,26,31-Heptaoxacyclo...	504	C28H56O7	066055-34-3	64
5	Methylamine, N-(1-methylhexylide...	127	C8H17N	022058-71-5	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

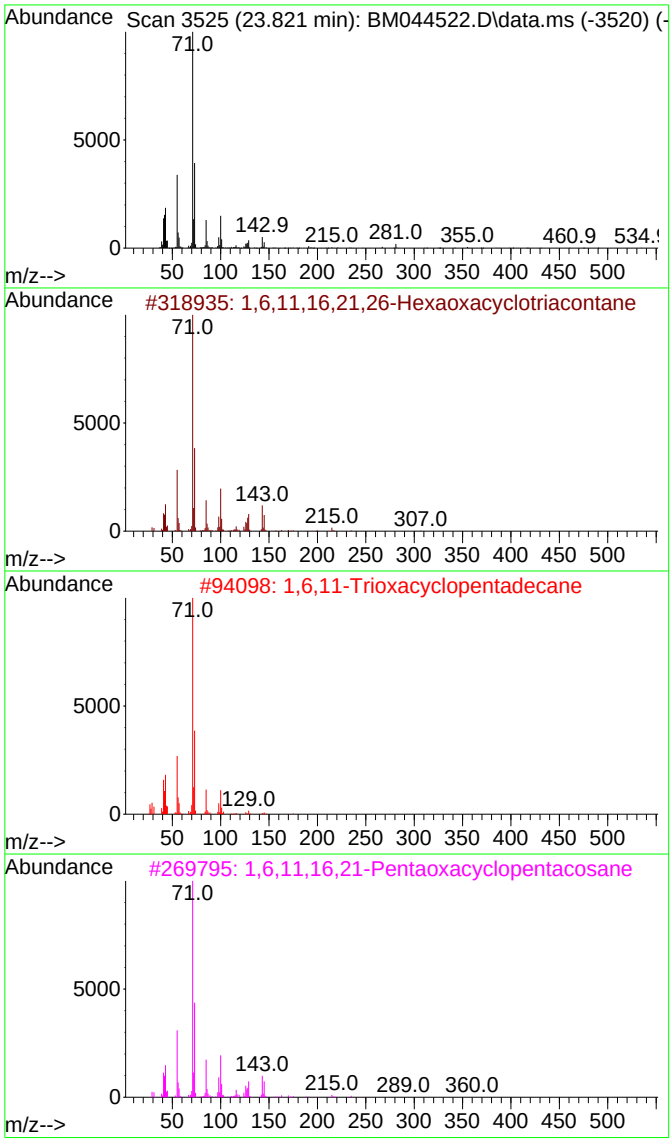
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

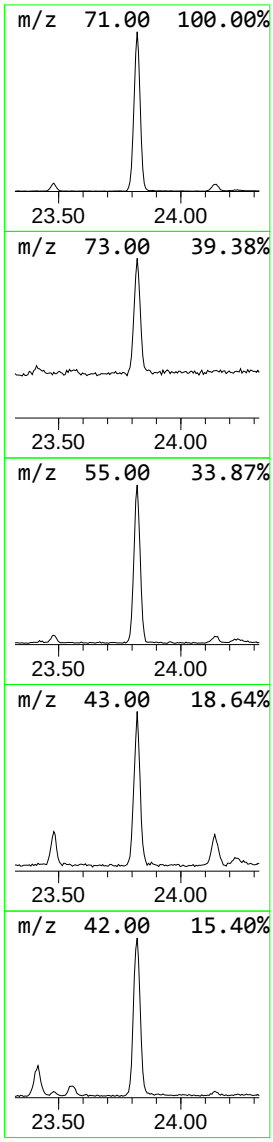
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 1,6,11,16,21,26-Hexaoxacycl... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.821	6.77 ng/ul	1019450	Perylene-d12	23.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,11,16,21,26-Hexaoxacyclotria...	432	C24H48O6	064001-05-4	91
2	1,6,11-Trioxacyclopentadecane	216	C12H24O3	000295-63-6	86
3	1,6,11,16,21-Pentaoxacyclopentac...	360	C20H40O5	056890-57-4	59
4	Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	58
5	1,6,11,16,21,26,31-Heptaoxacyclo...	504	C28H56O7	066055-34-3	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030624\
Data File : BM044522.D
Acq On : 06 Mar 2024 18:57
Operator : MA/JU
Sample : P1642-07
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0R35

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-brom...	9.222	15.0	ng/ul	1411440	2	10.528	1879780	20.0
Benzene, 1,2-di...	9.439	3.5	ng/ul	334041	2	10.528	1879780	20.0
Benzene, 1-chlo...	9.487	3.8	ng/ul	357451	2	10.528	1879780	20.0
Benzene, 1-brom...	9.528	8.1	ng/ul	762889	2	10.528	1879780	20.0
Benzene, 1-chlo...	11.122	28.9	ng/ul	2713880	2	10.528	1879780	20.0
Benzene, 1-chlo...	11.334	7.1	ng/ul	665577	2	10.528	1879780	20.0
4-Chloro-2-nitr...	11.810	2.2	ng/ul	205730	2	10.528	1879780	20.0
unknown-01	12.045	6.8	ng/ul	638491	2	10.528	1879780	20.0
unknown-02	18.492	8.1	ng/ul	1174210	4	17.115	2890180	20.0
1,6,11-Trioxacy...	18.986	2.0	ng/ul	291949	4	17.115	2890180	20.0
1,6,11,16,21-Pe...	21.409	5.2	ng/ul	782453	5	21.303	3033840	20.0
1,6,11,16,21,26...	23.821	6.8	ng/ul	1019450	6	23.556	3011160	20.0