LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.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	3.091	38	43	45	rBV2	4358053	6101493	100.00%	17.417%
2	3.114	45	47	66	rVV2	4085334	5886221	96.47%	16.803%
3	3.249	66	70	79	rVV	41216	83293	1.37%	0.238%
4	3.320	79	82	87	rVV	22710	32134	0.53%	0.092%
5	3.367	87	90	101	rVB	10604	14775	0.24%	0.042%
6	3.596	125	129	134	rVB	4634	6373	0.10%	0.018%
7	5.317	418	422	431	rVB	10472	15854	0.26%	0.045%
8	7.403	770	777	790	rVB2	93434	255962	4.20%	0.731%
9	7.579	800	807	814	rVV	82639	135313	2.22%	0.386%
10	7.673	815	823	830	rVB	152142	245223	4.02%	0.700%
11	7.790	836	843	852	rVB	99107	164209	2.69%	0.469%
12	8.260	916	923	931	rBV	62540	107037	1.75%	0.306%
13	8.701	990	998	1007	rVB	250141	419852	6.88%	1.199%
14	8.965	1036	1043	1048	rBV	131277	228571	3.75%	0.652%
15	9.024	1048	1053	1057	rVV	264081	449201	7.36%	1.282%
16	9.071	1057	1061	1069	rVB	166370	285751	4.68%	0.816%
17	9.359	1103	1110	1116	rBV	110410	186390	3.05%	0.532%
18	9.435	1116	1123	1131	rVB	87697	155414	2.55%	0.444%
19	9.999	1210	1219	1226	rBV	179151	306975	5.03%	0.876%
20	10.158	1239	1246	1253	rBV	72204	127940	2.10%	0.365%
21	10.240	1253	1260	1267	rBV	219178	379191	6.21%	1.082%
22	10.481	1294	1301	1311	rVB	177793	302725	4.96%	0.864%
23	10.704	1332	1339	1350	rBV2	136946	345556	5.66%	0.986%
24	11.092	1397	1405	1409	rBV	107020	198998	3.26%	0.568%
25	11.139	1409	1413	1420	rVV	154968	265135	4.35%	0.757%
26	11.215	1420	1426	1435	rVB	34573	59677	0.98%	0.170%
27	11.350	1445	1449	1454	rBV	8112	12614	0.21%	0.036%
28	11.421	1454	1461	1470	rVB	231419	390872	6.41%	1.116%
29	11.779	1516	1522	1527	rBV	10411	15842	0.26%	0.045%
30	11.844	1527	1533	1538	rVB	22207	33246	0.54%	0.095%
31	11.991	1550	1558	1573	rBV	87680	222592	3.65%	0.635%
32	12.578	1651	1658	1667	rBV	255738	387903	6.36%	1.107%
33	12.713	1677	1681	1689	rVB	10115	14652	0.24%	0.042%
34	12.842	1698	1703	1709	rBV2	8841	12471	0.20%	0.036%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Title : SVOA CALIBRATION

35	12.942	1712	1720	1729	rBV	401982	643103	10.54%	1.836%
36	13.083	1735	1744	1751	rVB2	239911	574350	9.41%	1.640%
37	13.389	1788	1796	1805	rBV	226227	358124	5.87%	1.022%
38	13.759	1853	1859	1867	rVB	162026	254270	4.17%	0.726%
39	14.276	1940	1947	1951	rBV	260467	393914	6.46%	1.124%
40	14.317	1951	1954	1961	rVB	168749	231474	3.79%	0.661%
41	14.575	1991	1998	2001	rBV	330337	578912	9.49%	1.653%
42	14.605	2001	2003	2011	rVB	248031	286975	4.70%	0.819%
43	14.769	2024	2031	2039	rBV	41970	67925	1.11%	0.194%
44	14.875	2041	2049	2055	rBV2	182836	298741	4.90%	0.853%
45	14.940	2055	2060	2064	rVV	408925	649844	10.65%	1.855%
46	14.975	2064	2066	2074	rVV	94236	123371	2.02%	0.352%
47	15.051	2074	2079	2087	rVV	114448	183513	3.01%	0.524%
48	15.163	2090	2098	2107	rVV	474152	723602	11.86%	2.066%
49	15.268	2110	2116	2125	rVB	189871	290563	4.76%	0.829%
50	15.721	2186	2193	2198	rBV	12281	16947	0.28%	0.048%
51	15.862	2210	2217	2221	rBV	439630	660602	10.83%	1.886%
52	15.909	2221	2225	2228	rBV2	258761	445668	7.30%	1.272%
53	15.985	2233	2238	2250	rVV2	294514	519030	8.51%	1.482%
54	17.254	2448	2454	2464	rBV	472230	712654	11.68%	2.034%
55	17.612	2508	2515	2521	rBV2	234022	347936	5.70%	0.993%
56	17.712	2525	2532	2535	rVV2	453712	708994	11.62%	2.024%
57	17.747	2535	2538	2545	rVB	475628	638188	10.46%	1.822%
58	18.012	2576	2583	2590	rBV	467777	705758	11.57%	2.015%
59	18.546	2669	2674	2679	rVB2	17895	23641	0.39%	0.067%
60	18.764	2706	2711	2715	rBV	22807	28511	0.47%	0.081%
61	18.805	2715	2718	2723	rVB2	15106	16788	0.28%	0.048%
62	19.010	2748	2753	2759	rVB	20938	29444	0.48%	0.084%
63	19.645	2852	2861	2867	rBV	319975	441382	7.23%	1.260%
64	19.974	2911	2917	2926	rBV	491834	713677	11.70%	2.037%
65	20.873	3064	3070	3075	rBV	341572	401496	6.58%	1.146%
66	21.501	3173	3177	3185	rVB4	6816	11454	0.19%	0.033%
67	21.589	3188	3192	3197	rVB	8870	13607	0.22%	0.039%
68	21.754	3214	3220	3226	rBV	361288	470616	7.71%	1.343%
69	21.889	3234	3243	3247	rBV	221294	383167	6.28%	1.094%
70	21.942	3247	3252	3259	rVB	460657	760322	12.46%	2.170%
71	22.065	3269	3273	3279	rVB	7870	11641	0.19%	0.033%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Title : SVOA CALIBRATION

72	22.705	3378	3382	3387	rVB4	7720	12488	0.20%	0.036%
73	22.970	3420	3427	3433	rVB	33802	63427	1.04%	0.181%
74	23.434	3500	3506	3516	rVB3	9637	22423	0.37%	0.064%
75	24.262	3636	3647	3660	rVB	372800	934405	15.31%	2.667%
76	25.038	3767	3779	3785	rBV	249639	713756	11.70%	2.037%
77	25.114	3785	3792	3804	rVV	315444	866252	14.20%	2.473%
78	25.267	3808	3818	3831	rVB2	133518	383540	6.29%	1.095%
79	26.483	4017	4025	4037	rVB6	11267	34434	0.56%	0.098%
80	27.916	4264	4269	4285	rVB9	9873	34356	0.56%	0.098%
81	29.126	4461	4475	4492	rVB2	89891	396415	6.50%	1.132%

Sum of corrected areas: 35031155

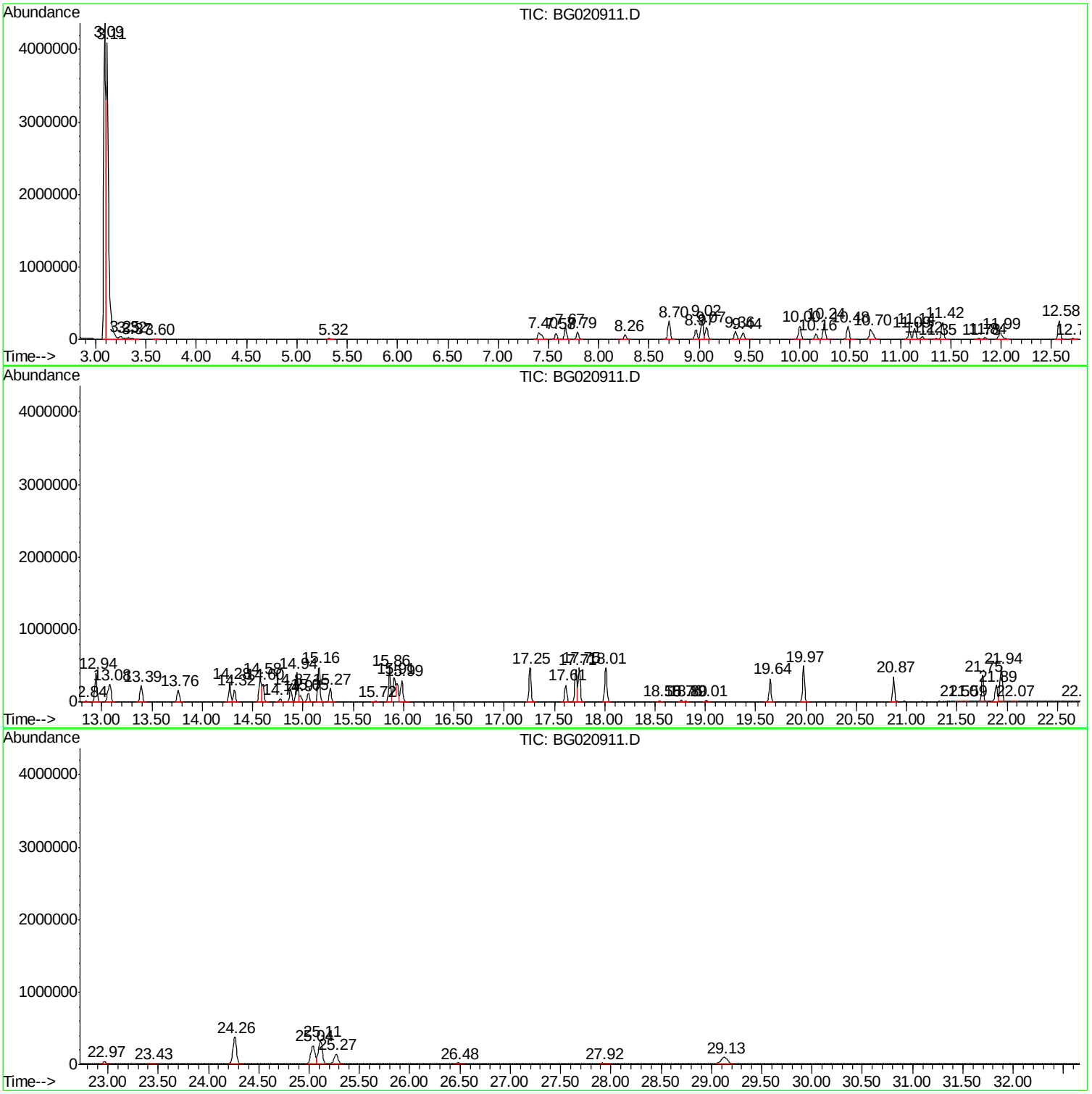
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

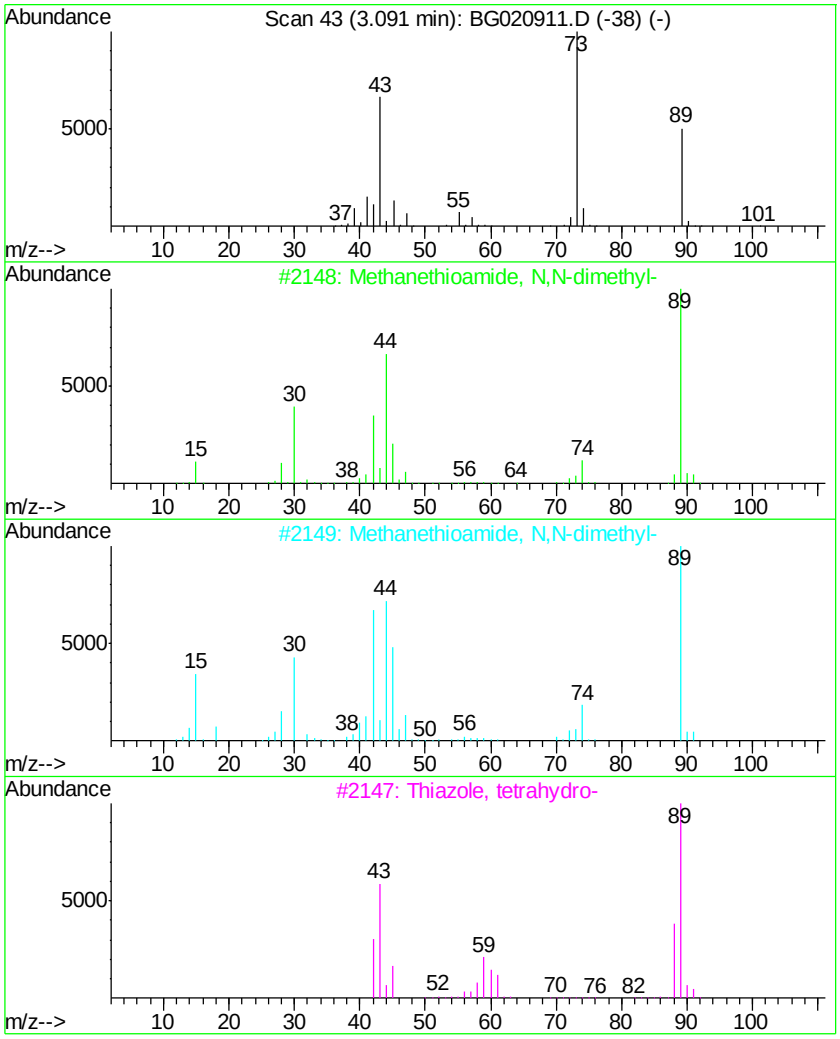
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

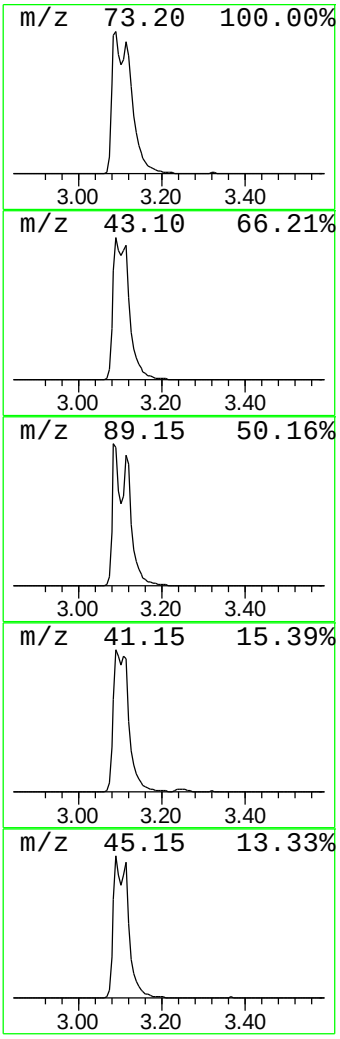
****

Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.09	1140.07 ng/ul	6101490	1,4-Dichlorobenzene-d4	8.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	47
2	Methanethioamide, N,N-dimethyl-	89	C3H7NS	000758-16-7	42
3	Thiazole, tetrahydro-	89	C3H7NS	000504-78-9	28
4	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	25
5	1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

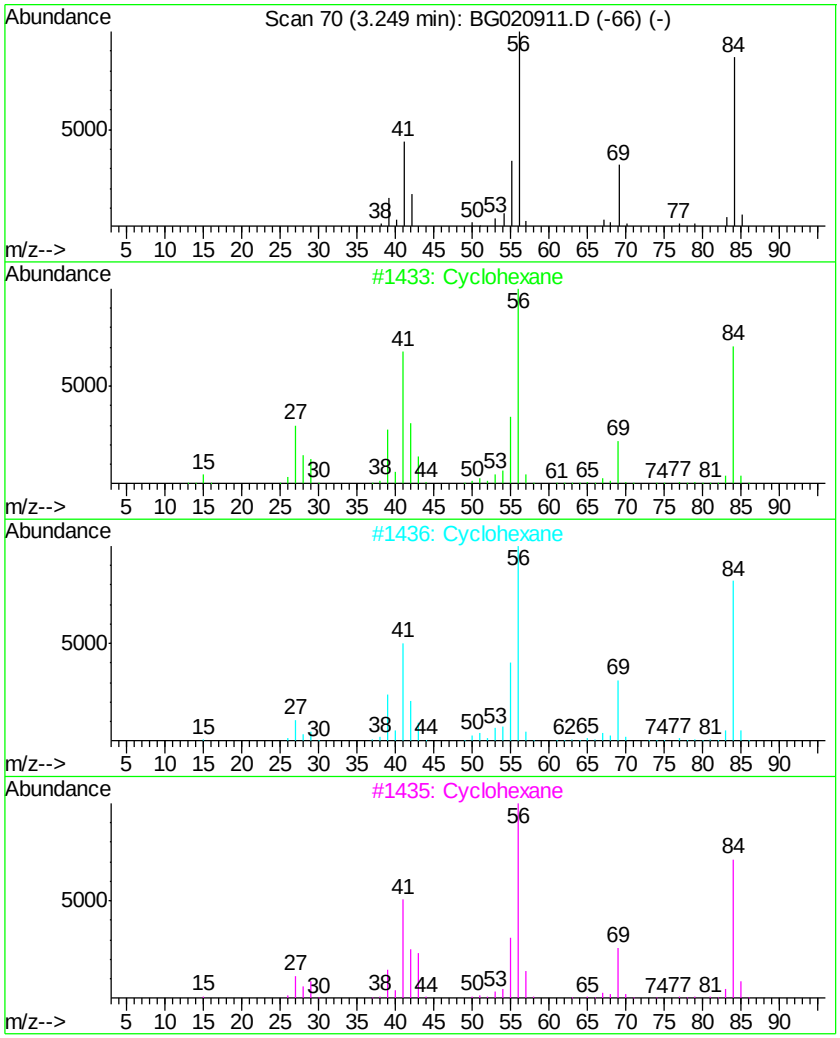
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

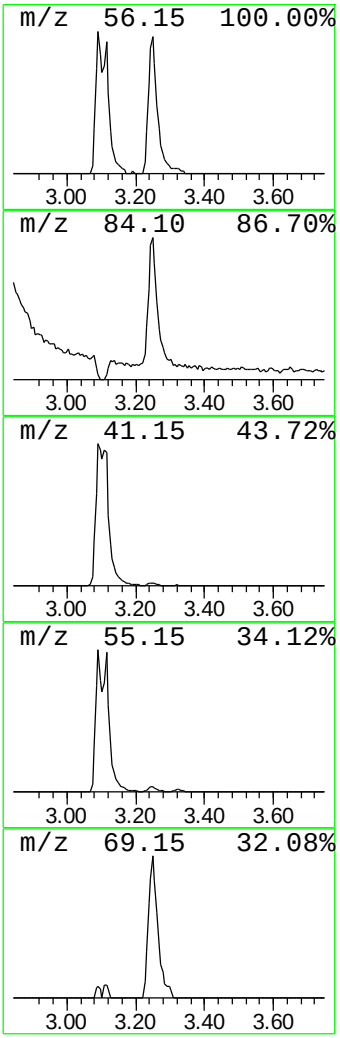
****

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.25	15.56 ng/ul	83293	1,4-Dichlorobenzene-d4	8.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane	84	C6H12	000110-82-7	91
2	Cyclohexane	84	C6H12	000110-82-7	91
3	Cyclohexane	84	C6H12	000110-82-7	91
4	Cyclohexane	84	C6H12	000110-82-7	64
5	2-Amino-oxazole	84	C3H4N2O	004570-45-0	53





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

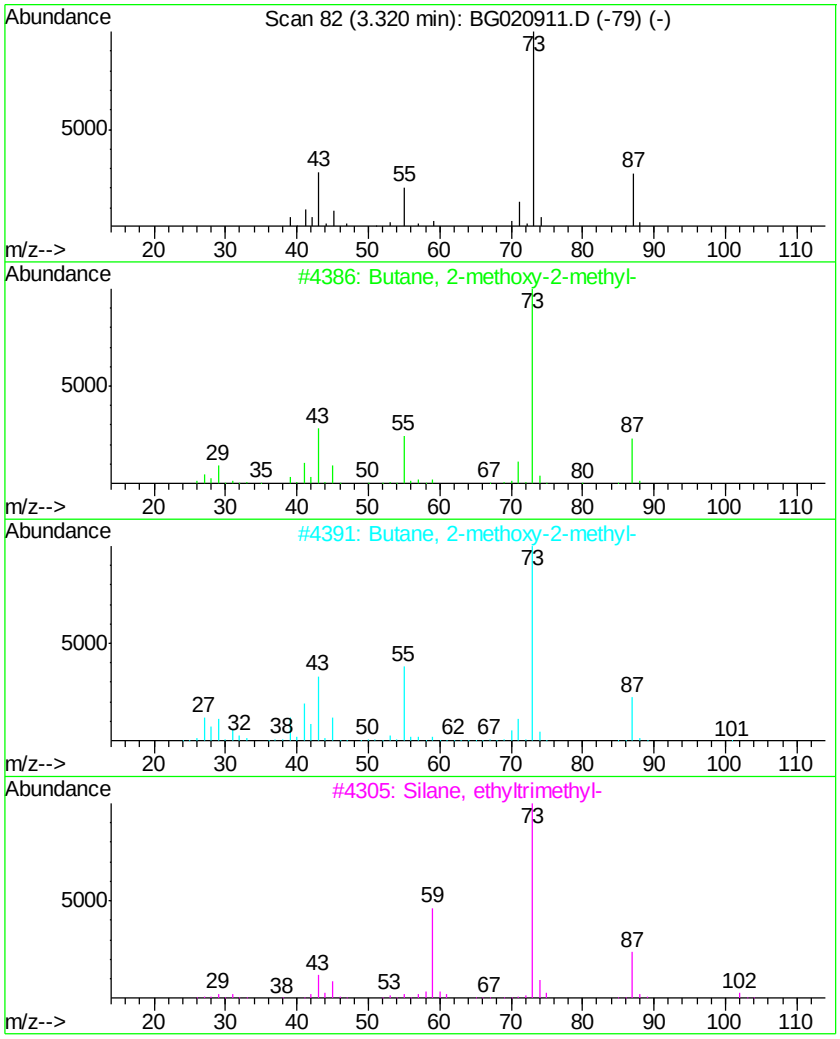
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

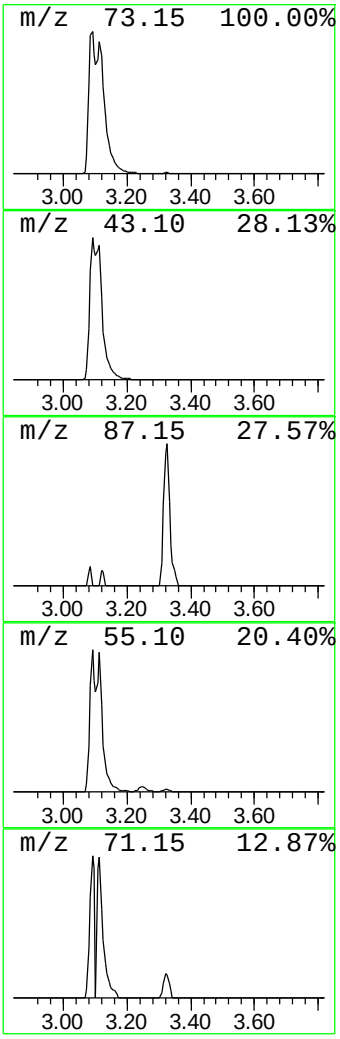
****

Peak Number 4 Butane, 2-methoxy-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.32	6.00 ng/ul	32134	1,4-Dichlorobenzene-d4	8.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	64
3	Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	36
4	2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	32
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

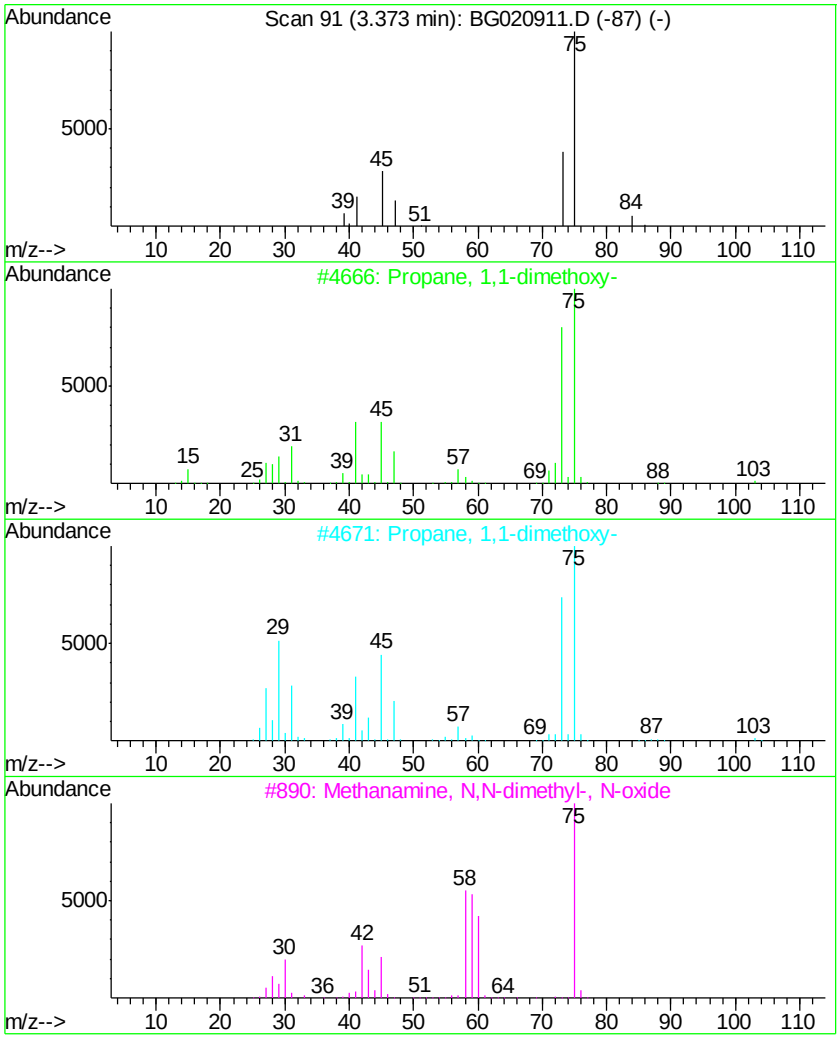
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

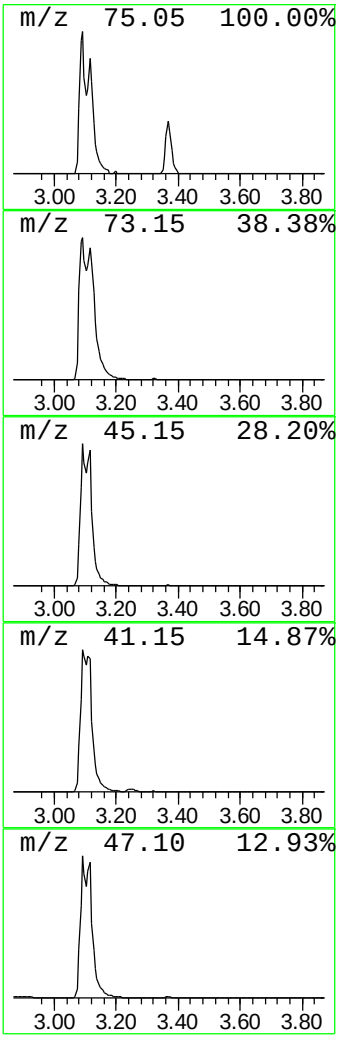
****

Peak Number 5 Propane, 1,1-dimethoxy- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.37	2.76 ng/ul	14775	1,4-Dichlorobenzene-d4	8.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	74
2	Propane, 1,1-dimethoxy-	104	C5H12O2	004744-10-9	9
3	Methanamine, N,N-dimethyl-, N-oxide	75	C3H9NO	001184-78-7	4
4	Formamide, N-methylthio	75	C2H5NS	018952-41-5	4
5	Glycine	75	C2H5NO2	000056-40-6	4





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

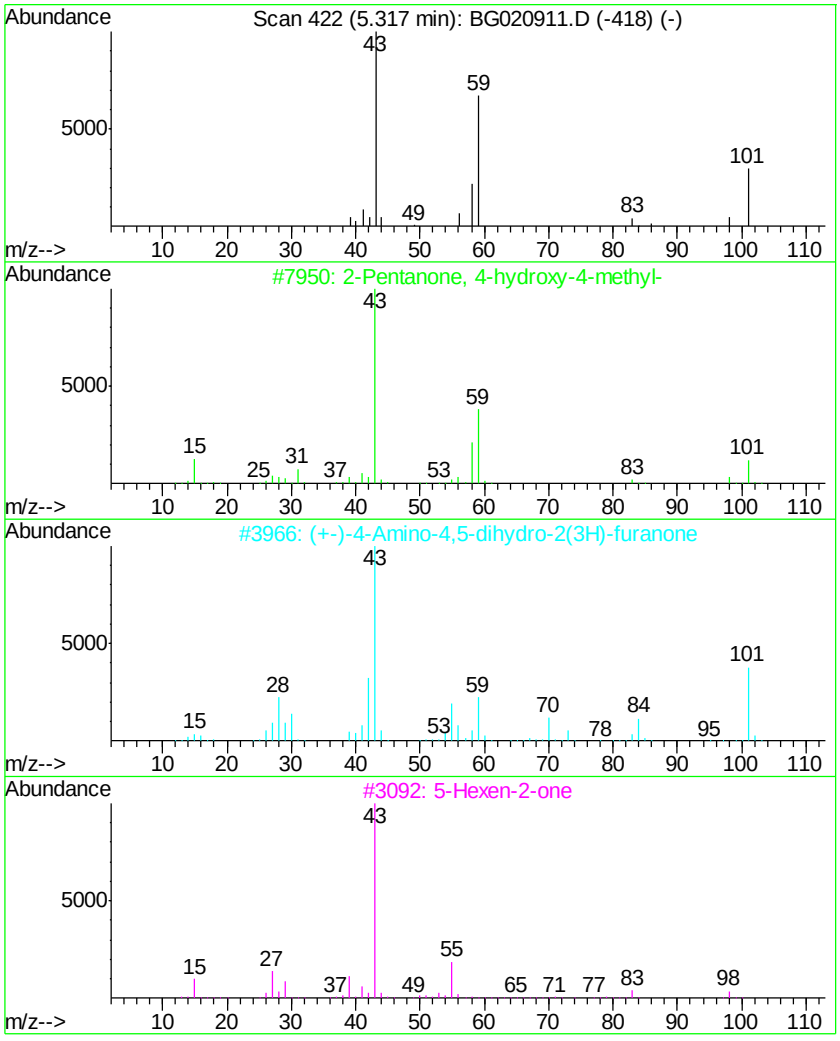
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

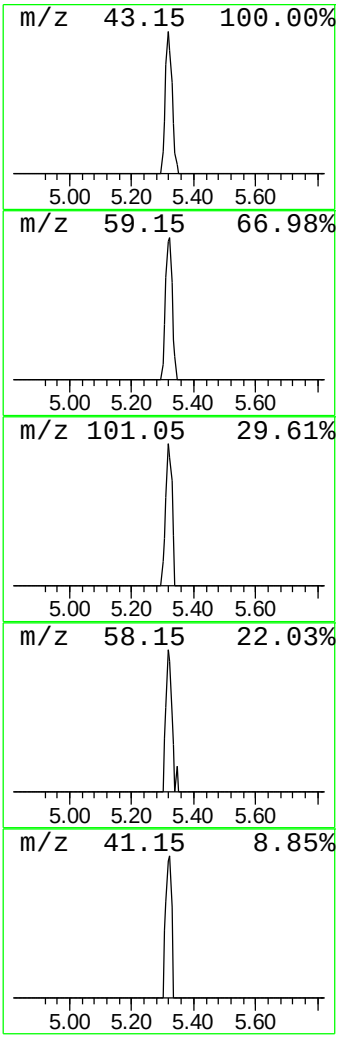
****

Peak Number 6 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.32	2.96 ng/ul	15854	1,4-Dichlorobenzene-d4	8.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	5-Hexen-2-one	98	C6H10O	000109-49-9	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

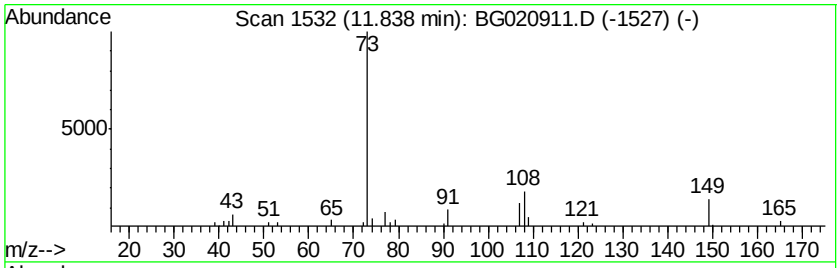
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

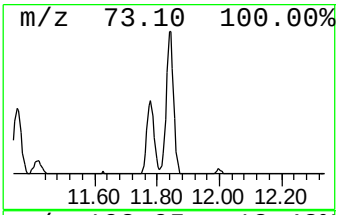
****

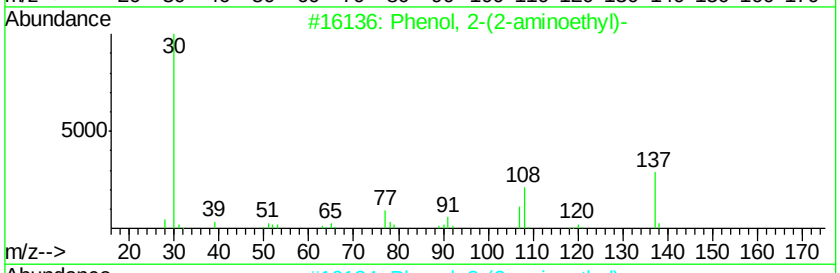
Peak Number 7 unknown-03 Concentration Rank 8

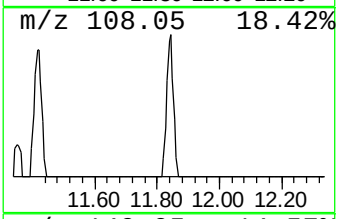
R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	3.34 ng/ul	33246	Naphthalene-d8	11.09

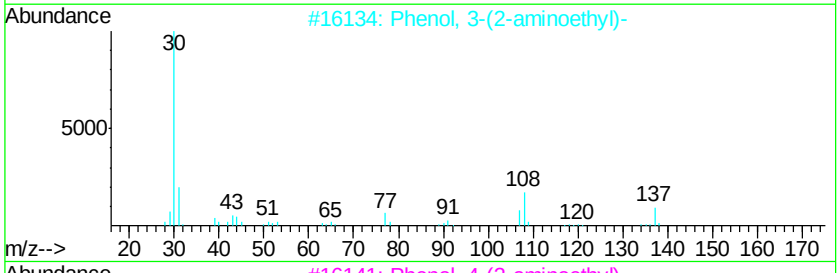
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2-(2-aminoethyl)-	137	C8H11NO	002039-66-9	16
2	Phenol, 3-(2-aminoethyl)-	137	C8H11NO	000588-05-6	9
3	Phenol, 4-(2-aminoethyl)-	137	C8H11NO	000051-67-2	9
4	1,3-Dioxolane, 2-(chloromethyl)-	122	C4H7ClO2	002568-30-1	9
5	N-Ethylformamide	73	C3H7NO	000627-45-2	7

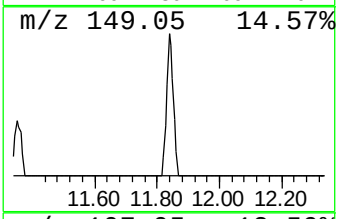


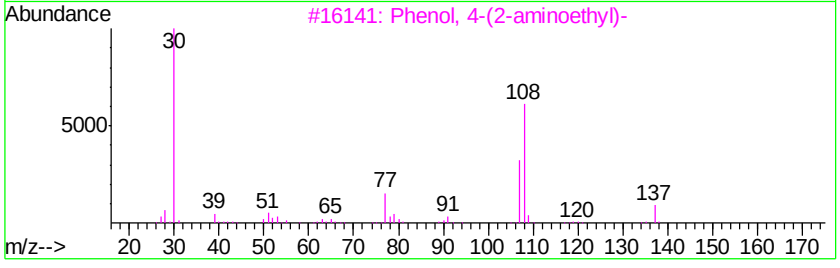


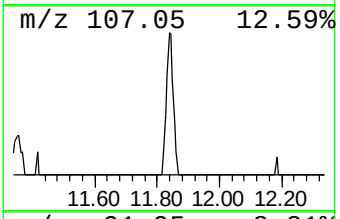


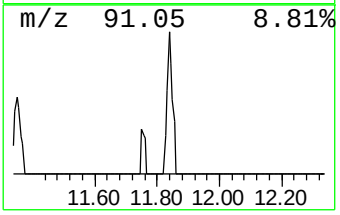












Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

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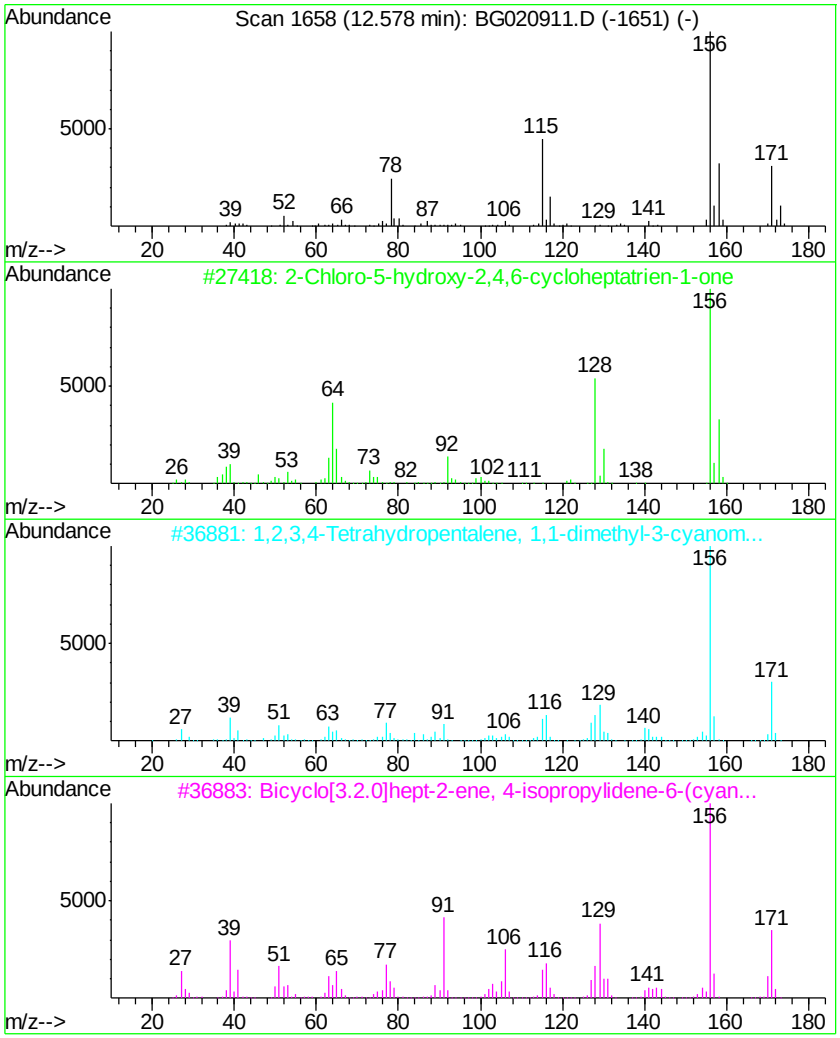
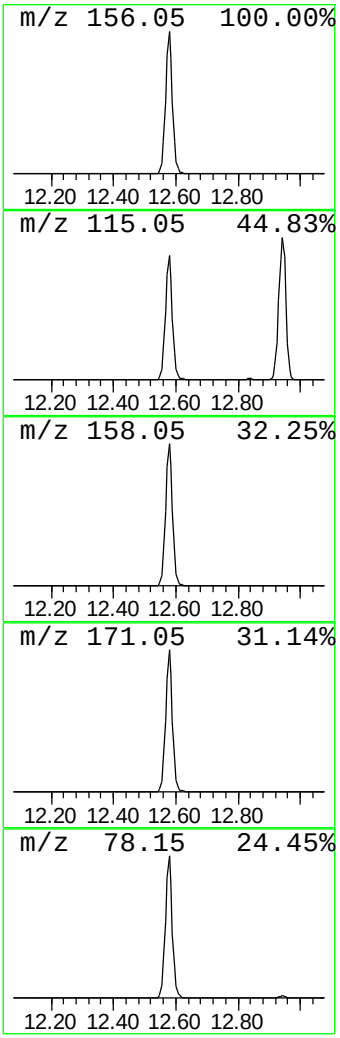
Peak Number 8 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.58	38.99 ng/ul	387903	Naphthalene-d8	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	38
2	1,2,3,4-Tetrahydropentalene, 1,1...	171	C12H13N	1000217-18-6	37
3	Bicyclo[3.2.0]hept-2-ene, 4-isop...	171	C12H13N	1000217-15-6	32
4	6-Isopropylquinoline	171	C12H13N	000135-79-5	32
5	Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	28





Library Search Compound Report

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Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

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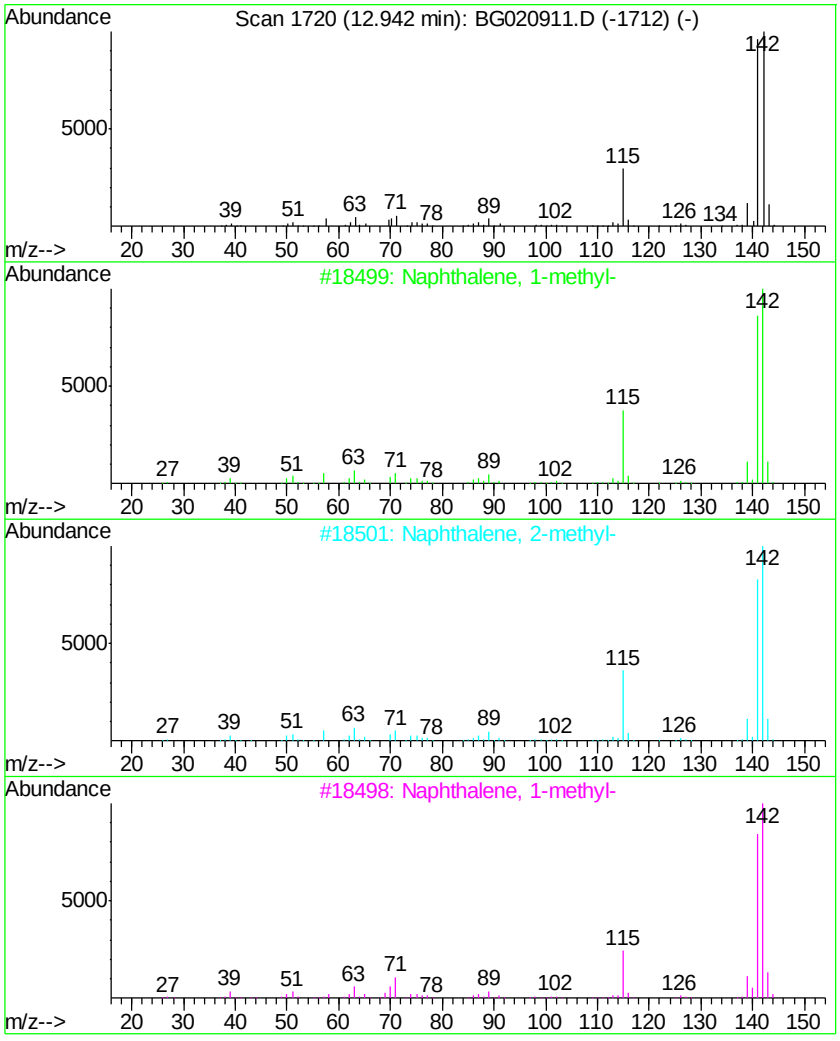
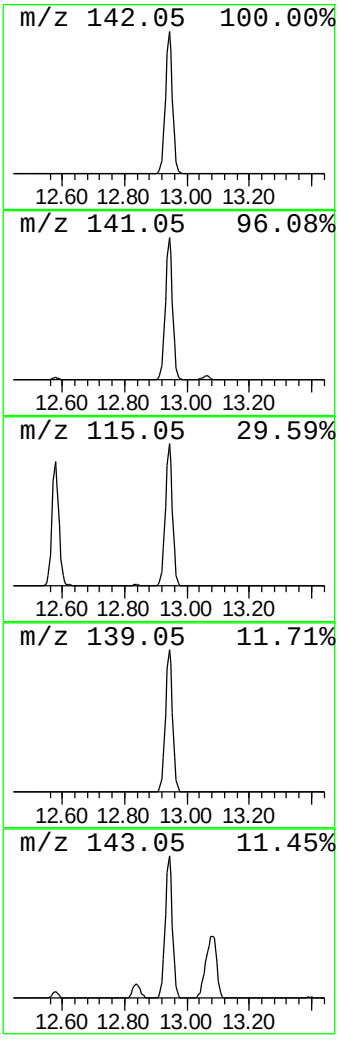
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.94	64.63 ng/ul	643103	Naphthalene-d8	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	Benzocycloheptatriene	142	C11H10	000264-09-5	91





Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

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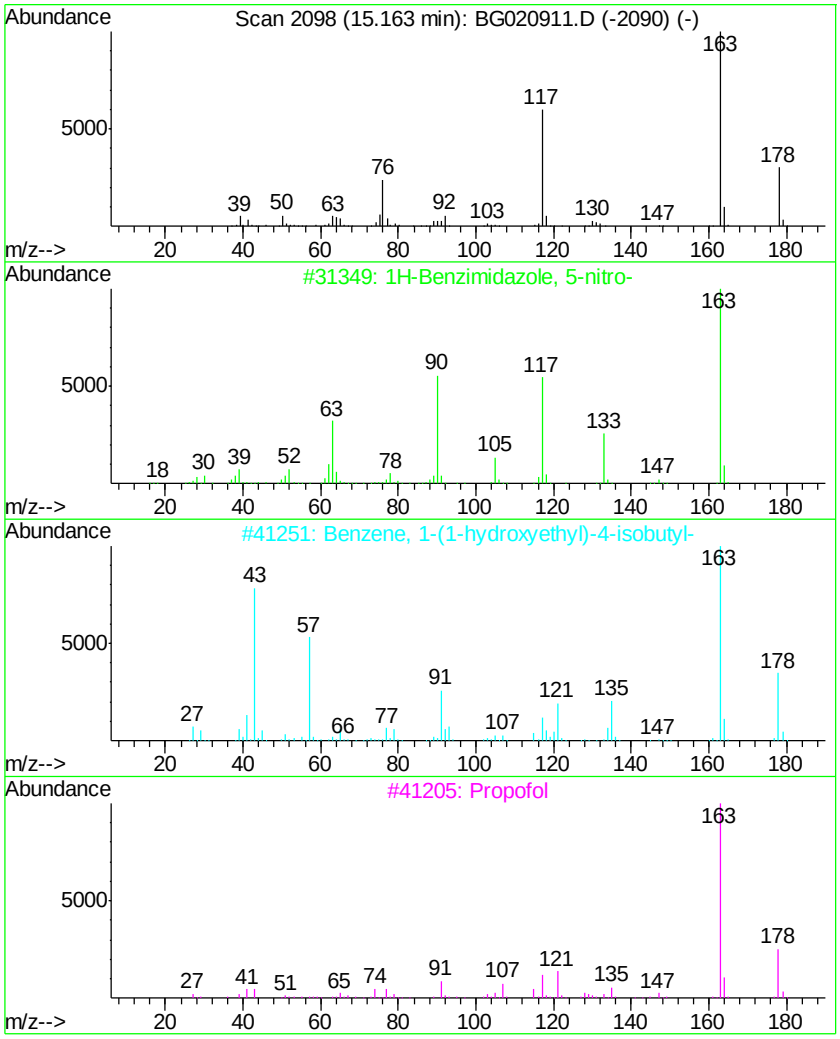
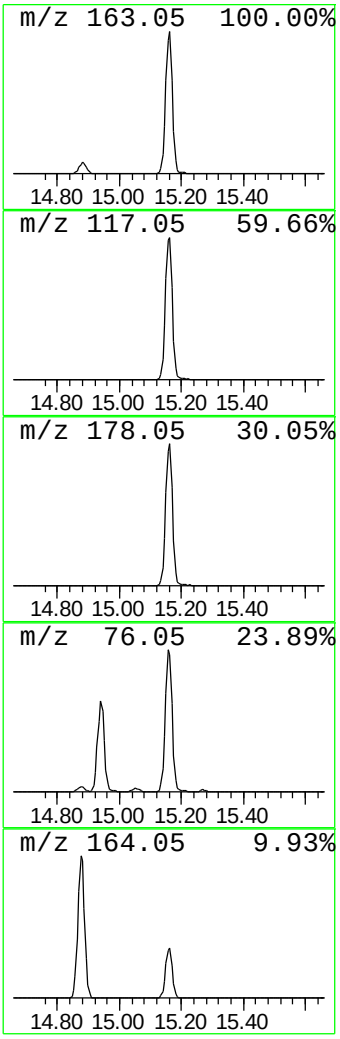
Peak Number 10 1H-Benzimidazole, 5-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	48.44 ng/ul	723602	Acenaphthene-d10	14.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 5-nitro-	163	C7H5N3O2	000094-52-0	53
2	Benzene, 1-(1-hydroxyethyl)-4-is...	178	C12H18O	040150-92-3	50
3	Propofol	178	C12H18O	002078-54-8	50
4	Ethanone, 1-(2,3-dihydro-1,4-ben...	178	C10H10O3	002879-20-1	49
5	Benzene, 1-(1,1-dimethylethyl)-2...	178	C12H18O	000088-40-4	47





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Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

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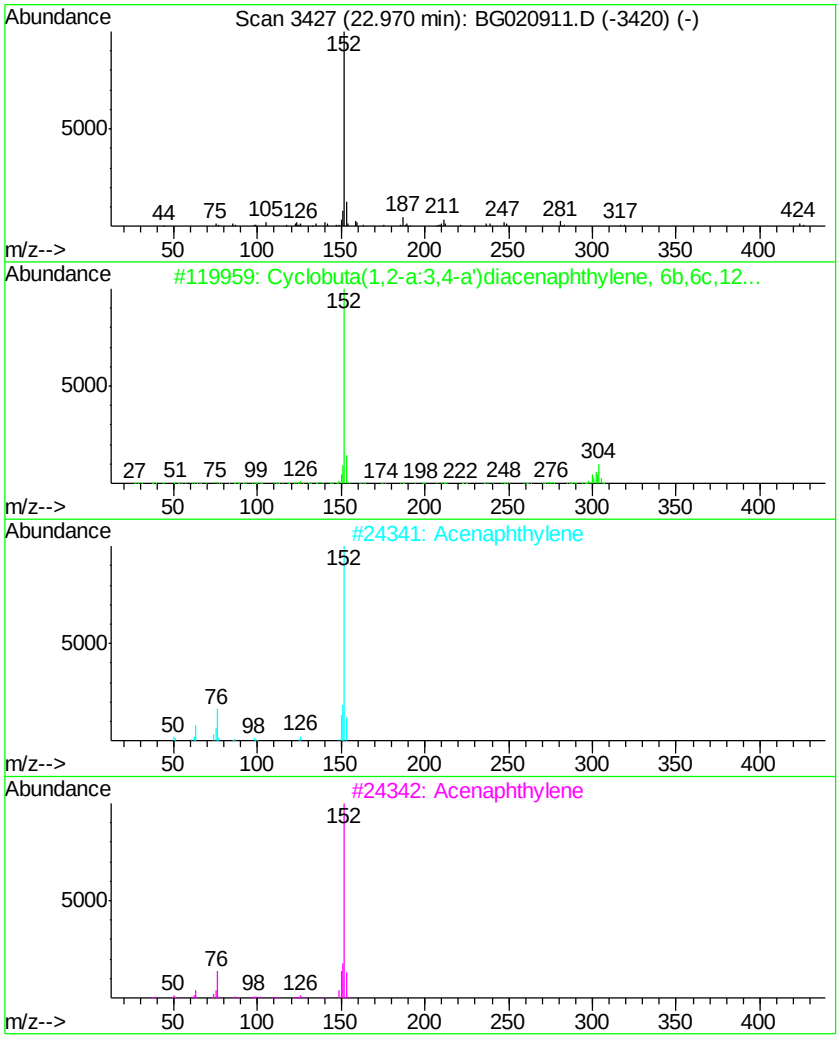
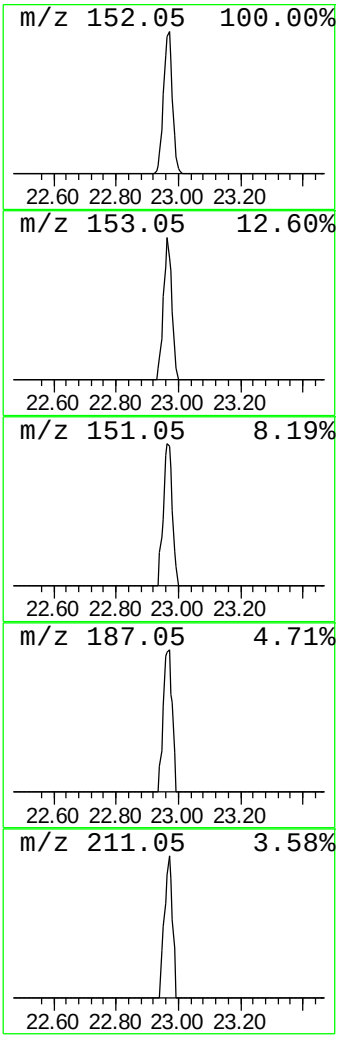
Peak Number 11 Cyclobuta(1,2-a:3,4-a')diac... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.97	3.31 ng/ul	63427	Chrysene-d12	21.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobuta(1,2-a:3,4-a')diacenaph...	304	C24H16	007259-03-2	50
2	Acenaphthylene	152	C12H8	000208-96-8	50
3	Acenaphthylene	152	C12H8	000208-96-8	39
4	[2-(Naphth-2-yl)vinyl]-methyl su...	232	C13H12O2S	1000286-42-3	39
5	6-Mercaptopurine	152	C5H4N4S	000050-44-2	9





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021616\
Data File : BG020911.D
Acq On : 16 Feb 2016 16:45
Operator : SJ/UM
Sample : H1454-06
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : D9QT5

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG020816.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown-01	3.09	1140.1	ng/ul	6101490	1	8.26	107037	20.0
(DEL) Alkane: Str...	3.25	15.6	ng/ul	83293	1	8.26	107037	20.0
Butane, 2-methoxy...	3.32	6.0	ng/ul	32134	1	8.26	107037	20.0
Propane, 1,1-dime...	3.37	2.8	ng/ul	14775	1	8.26	107037	20.0
unknown-02	5.32	3.0	ng/ul	15854	1	8.26	107037	20.0
unknown-03	11.84	3.3	ng/ul	33246	2	11.09	198998	20.0
unknown-04	12.58	39.0	ng/ul	387903	2	11.09	198998	20.0
Naphthalene, 1-me...	12.94	64.6	ng/ul	643103	2	11.09	198998	20.0
1H-Benzimidazole,...	15.16	48.4	ng/ul	723602	3	14.87	298741	20.0
Cyclobuta(1,2-a:3...	22.97	3.3	ng/ul	63427	5	21.89	383167	20.0