

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
 Data File : BM004545.D
 Acq On : 03 Mar 2016 23:13
 Operator : SJ/UM
 Sample : H1616-06
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BLDG06-P003-SS001-01

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.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	6.834	700	705	716	rBV	61213	112798	3.20%	0.427%
2	6.934	716	722	732	rVB	61812	93620	2.65%	0.354%
3	7.151	753	759	768	rBB	85271	131839	3.74%	0.499%
4	7.592	828	834	841	rBB	106689	162898	4.62%	0.617%
5	8.357	959	964	977	rBB	92493	159720	4.53%	0.605%
6	8.757	1026	1032	1041	rBB	77413	126044	3.57%	0.477%
7	9.475	1149	1154	1162	rBB	69738	110483	3.13%	0.418%
8	10.063	1249	1254	1269	rBV	92529	182054	5.16%	0.689%
9	10.380	1301	1308	1313	rBV	178812	292147	8.28%	1.106%
10	10.428	1313	1316	1323	rVB	30264	45475	1.29%	0.172%
11	10.533	1329	1334	1350	rBV	63670	119592	3.39%	0.453%
12	13.674	1862	1868	1872	rBV	248344	336631	9.55%	1.274%
13	13.721	1872	1876	1882	rVB	91218	123219	3.49%	0.466%
14	13.951	1909	1915	1927	rBB	292014	431420	12.23%	1.633%
15	14.263	1961	1968	1974	rVB2	282656	440542	12.49%	1.668%
16	15.110	2106	2112	2116	rVB	35508	41312	1.17%	0.156%
17	15.257	2131	2137	2144	rBV	379419	552621	15.67%	2.092%
18	15.415	2159	2164	2171	rBV	43360	69135	1.96%	0.262%
19	15.974	2254	2259	2261	rBV	25575	36798	1.04%	0.139%
20	17.015	2430	2436	2440	rBV	379576	520574	14.76%	1.971%
21	17.057	2440	2443	2447	rVV	51972	66992	1.90%	0.254%
22	17.115	2447	2453	2464	rVB	415944	602191	17.08%	2.280%
23	17.945	2591	2594	2603	rVB2	26049	38622	1.10%	0.146%
24	18.856	2746	2749	2753	rVB	35762	44473	1.26%	0.168%
25	19.086	2785	2788	2796	rVB	87266	125823	3.57%	0.476%
26	19.427	2841	2846	2849	rBV	424642	607480	17.23%	2.300%
27	21.233	3147	3153	3161	rBV3	503271	1488940	42.22%	5.636%
28	21.492	3191	3197	3201	rBV	649835	1288855	36.55%	4.879%
29	21.527	3201	3203	3207	rVB	698379	786053	22.29%	2.976%
30	21.597	3211	3215	3225	rVB	831766	1120728	31.78%	4.242%
31	21.703	3226	3233	3236	rBV	1574217	3526242	100.00%	13.348%
32	21.733	3236	3238	3243	rVB2	1417646	1734027	49.17%	6.564%
33	21.792	3243	3248	3260	rVB	1330457	1777879	50.42%	6.730%
34	21.897	3261	3266	3268	rBV	984805	1315897	37.32%	4.981%

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35	21.927	3268	3271	3277	rVB	1094204	1274541	36.14%	4.825%
36	22.074	3291	3296	3300	rVB	959712	1233617	34.98%	4.670%
37	22.274	3324	3330	3341	rVB	1646360	2776497	78.74%	10.510%
38	22.374	3342	3347	3353	rBV	1331419	2101626	59.60%	7.956%
39	23.309	3501	3506	3510	rBV	219988	417687	11.85%	1.581%

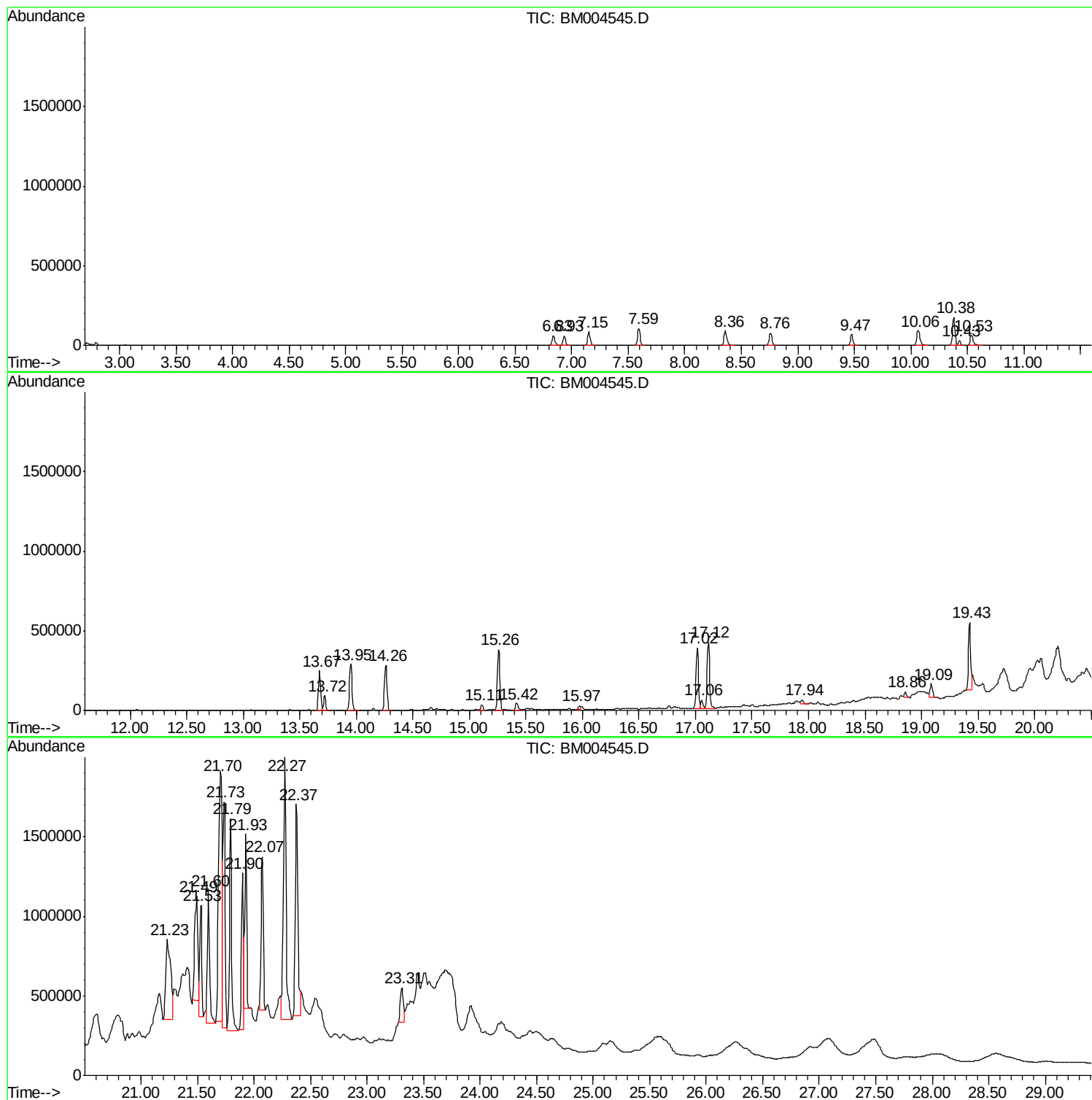
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Quant Title : SVOA CALIBRATION

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



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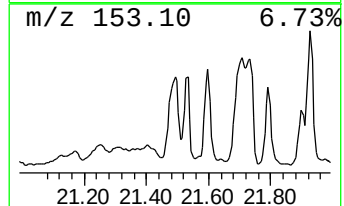
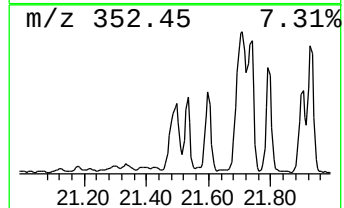
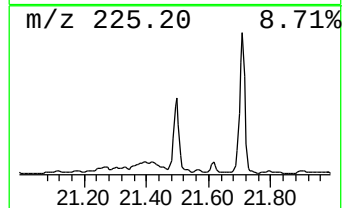
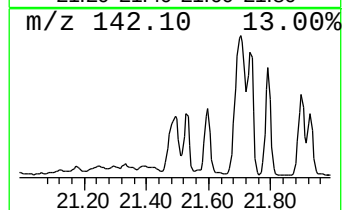
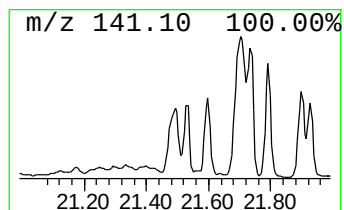
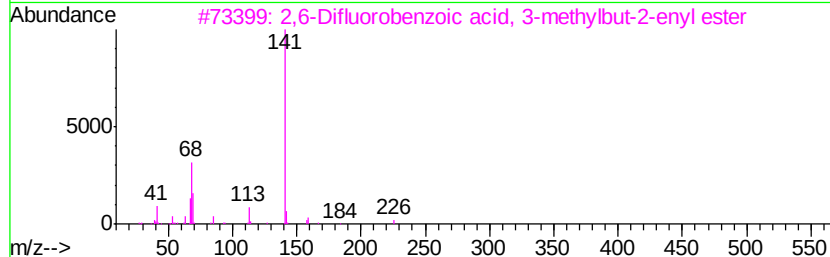
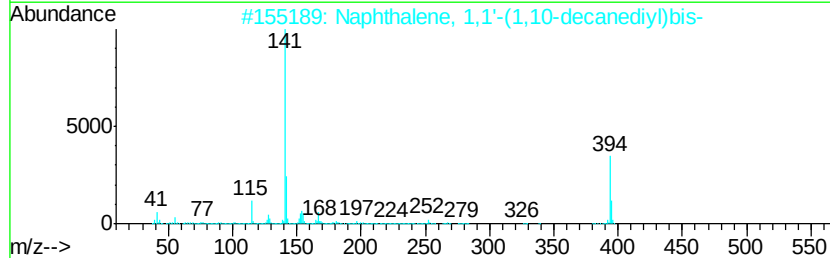
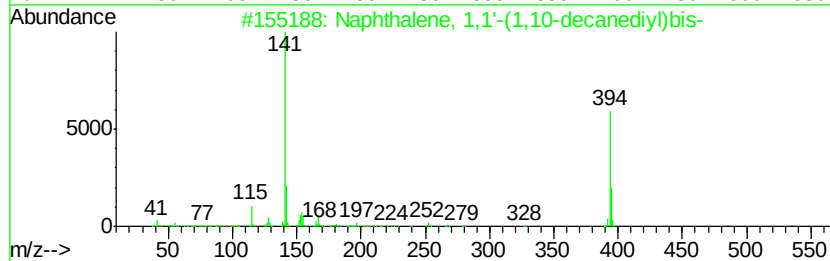
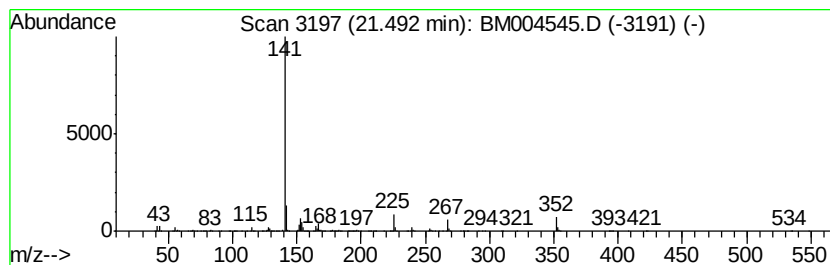
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Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.49	17.31 ng/ul	1288860	Chrysene-d12	21.23		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,1'-(1,10-decanedi...	394	C30H34	040339-27-3	35
2		Naphthalene, 1,1'-(1,10-decanedi...	394	C30H34	040339-27-3	17
3		2,6-Difluorobenzoic acid, 3-meth...	226	C12H12F2O2	1000292-58-2	9
4		1-Naphthalenepropionic acid	200	C13H12O2	003243-42-3	9
5		Pyrazol-4-amine, 3,5-dimethyl-1-...	251	C16H17N3	1000273-77-6	9



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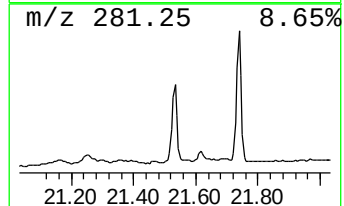
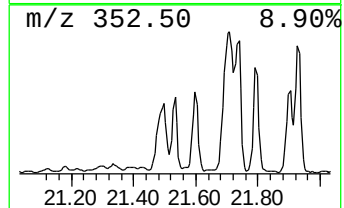
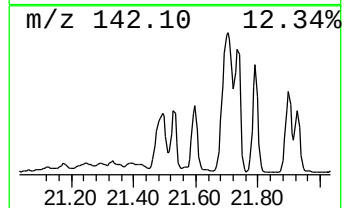
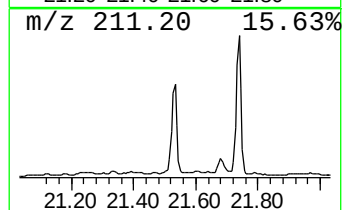
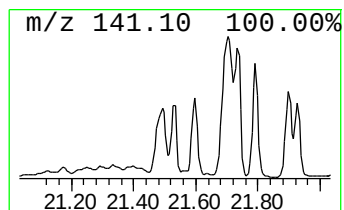
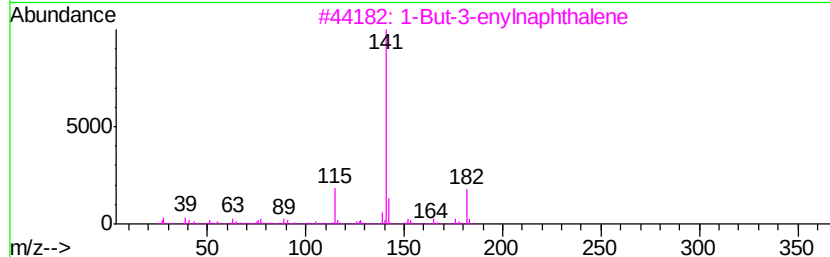
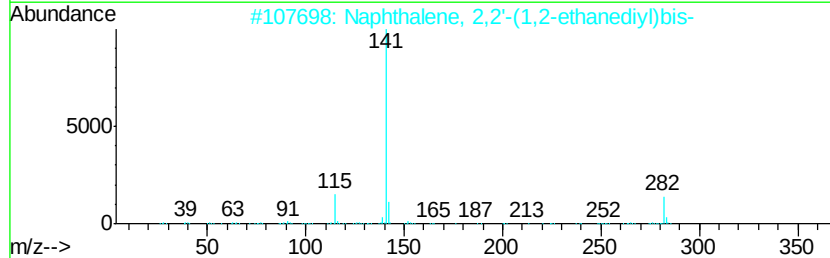
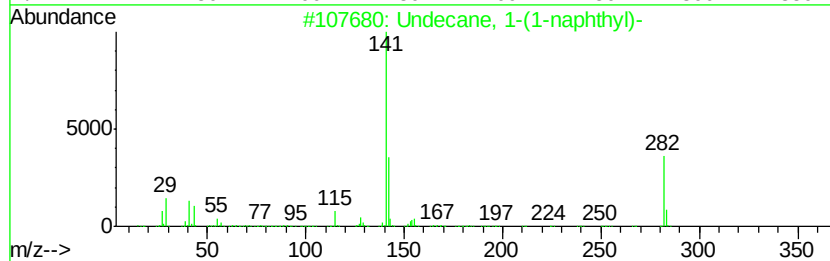
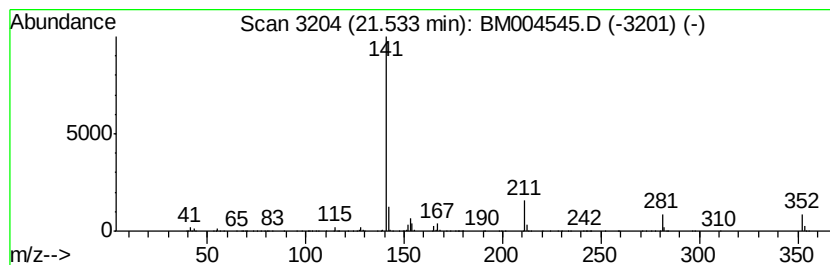
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TIC Library : C:\DATABASE\NIST02.L
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Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.53	10.56 ng/ul	786053	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Undecane, 1-(1-naphthyl)-	282	C21H30	007225-71-0	33
2		Naphthalene, 2,2'-(1,2-ethanediyl)bis-	282	C22H18	021969-45-9	9
3		1-But-3-enylnaphthalene	182	C14H14	002489-88-5	9
4		Naphthalene, 1-propyl-	170	C13H14	002765-18-6	9
5		Naphthalene, 2-hexyl-	212	C16H20	002876-46-2	9



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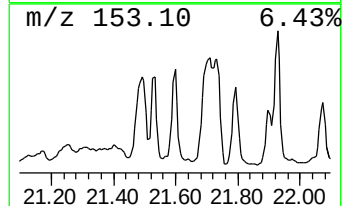
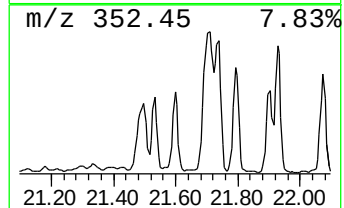
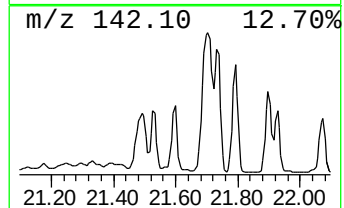
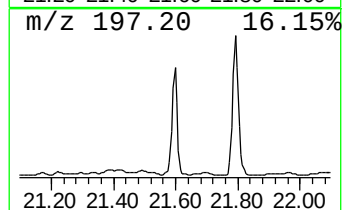
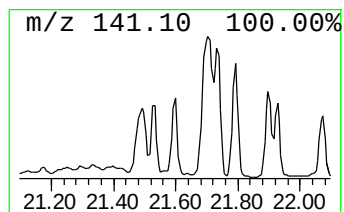
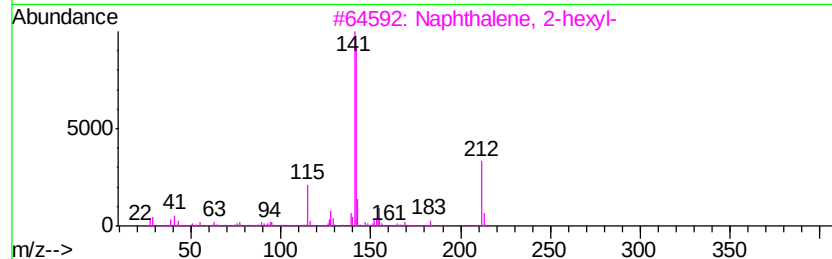
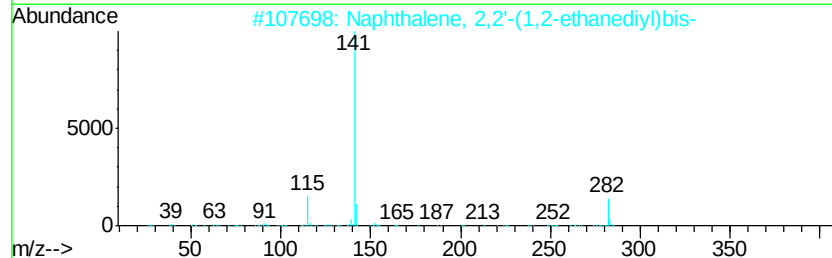
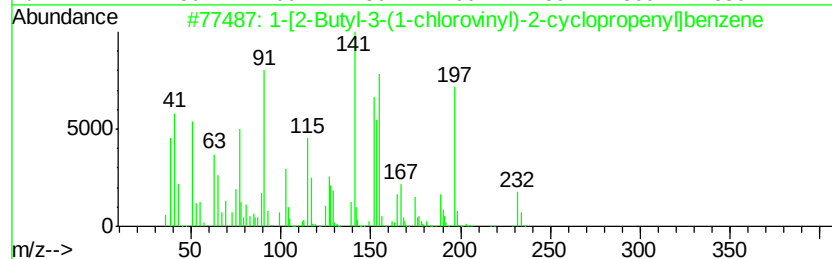
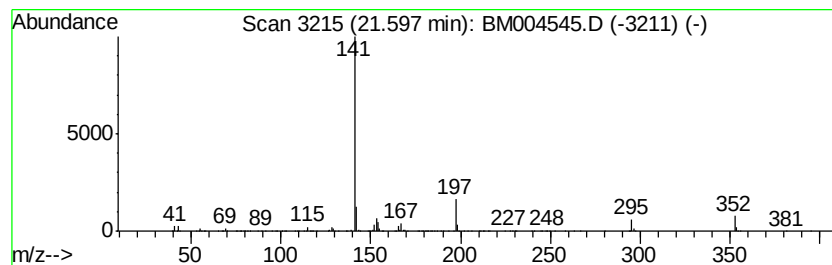
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Peak Number 3 unknown-02 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.60	15.05 ng/ul	1120730	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-[2-Butyl-3-(1-chlorovinyl)-2-c...	232	C15H17Cl	1000261-30-1	45
2		Naphthalene, 2,2'-(1,2-ethanediyl...	282	C22H18	021969-45-9	9
3		Naphthalene, 2-hexyl-	212	C16H20	002876-46-2	9
4		Naphthalene, 1-propyl-	170	C13H14	002765-18-6	9
5		1-But-3-enyl naphthalene	182	C14H14	002489-88-5	9



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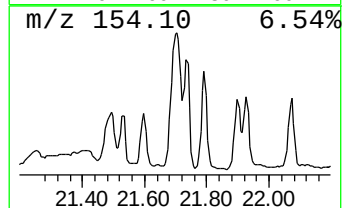
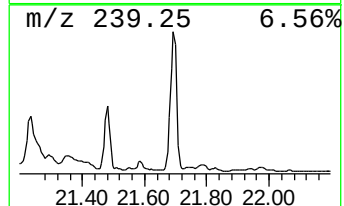
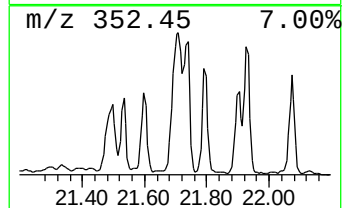
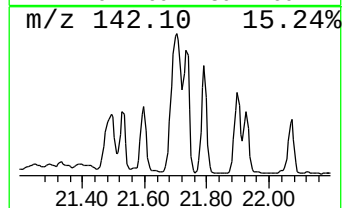
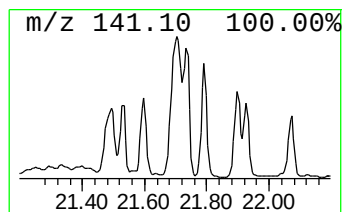
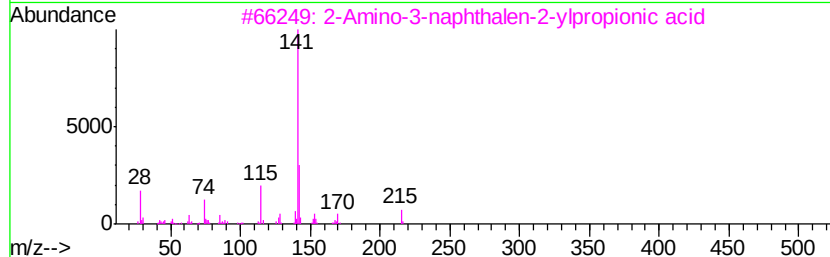
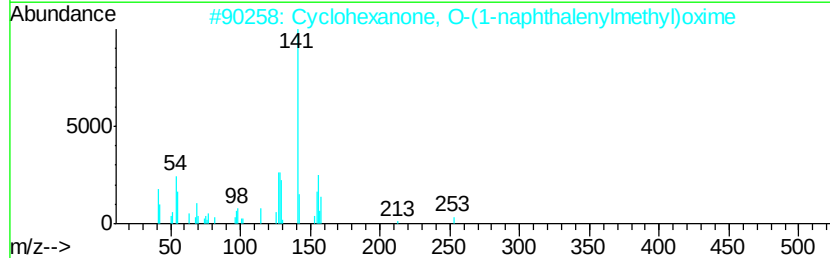
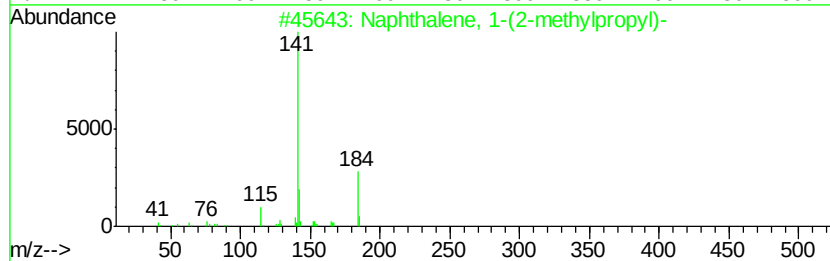
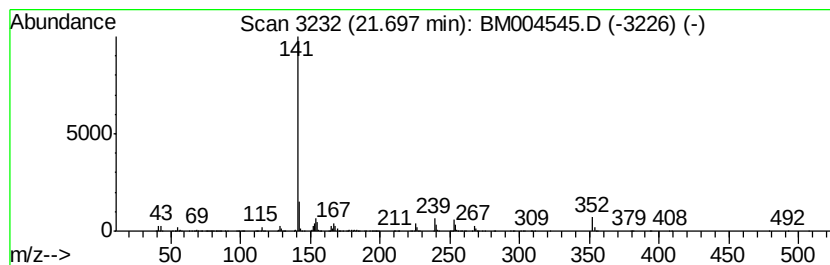
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Peak Number 4 Naphthalene, 1-(2-methylpro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.70	47.37 ng/ul	3526240	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1-(2-methylpropyl)-	184 C14H16	016727-91-6 59
2		Cyclohexanone, O-(1-naphthalenyl)-	253 C17H19NO	055045-02-8 9
3		2-Amino-3-naphthalen-2-ylpropion...	215 C13H13NO2	099631-78-4 9
4		Naphthalene, 2,2'-(1,2-ethanedi...	282 C22H18	021969-45-9 9
5		1-But-3-enyl naphthalene	182 C14H14	002489-88-5 9



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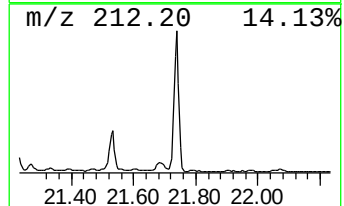
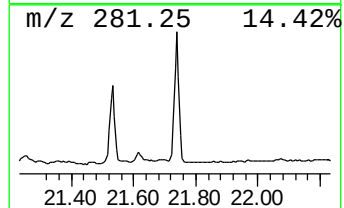
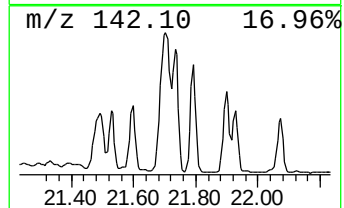
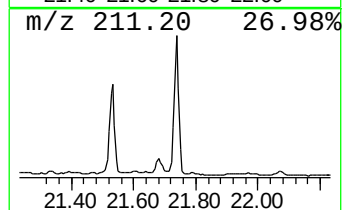
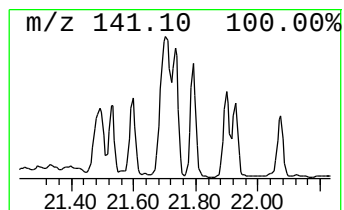
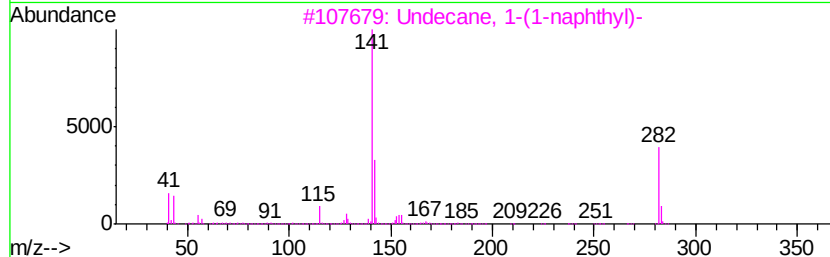
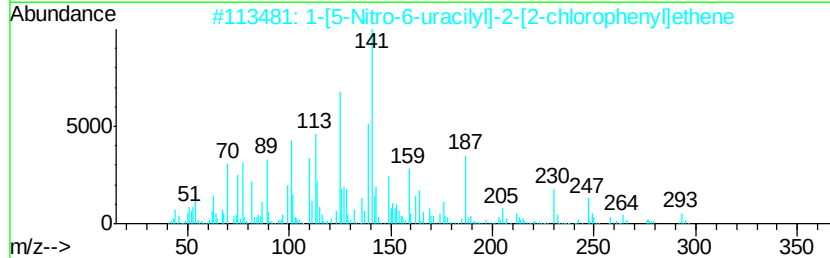
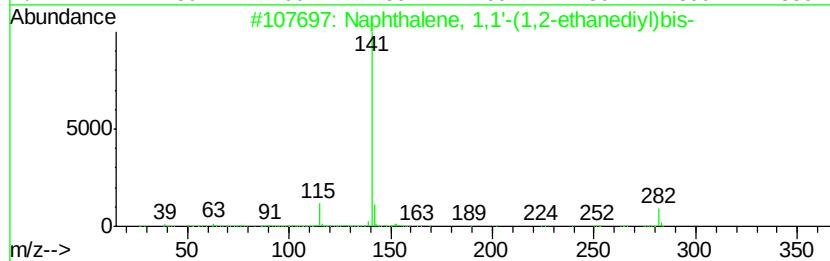
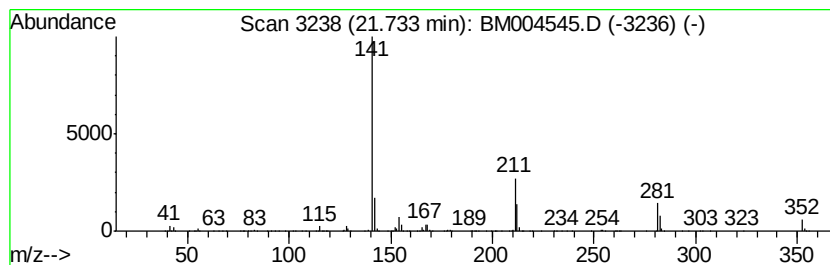
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Peak Number 5 unknown-03 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.73	23.29 ng/ul	1734030	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,1'-(1,2-ethanediyl)-	282	C22H18	015374-45-5	47
2		1-[5-Nitro-6-uracilyl]-2-[2-chlo...	293	C12H8ClN3O4	296798-53-3	38
3		Undecane, 1-(1-naphthyl)-	282	C21H30	007225-71-0	12
4		Naphthalene, 1-hexyl-	212	C16H20	002876-53-1	11
5		Pyrazol-4-amine, 3,5-dimethyl-1-...	251	C16H17N3	1000273-77-6	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BLDG06-P003-SS001-01

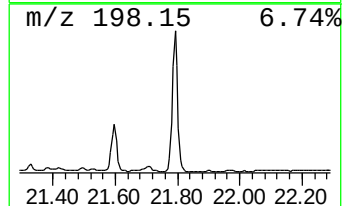
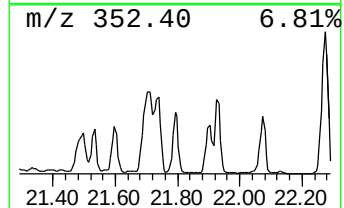
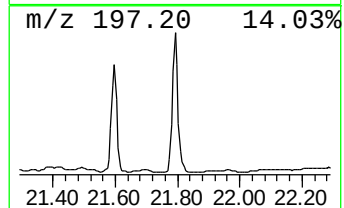
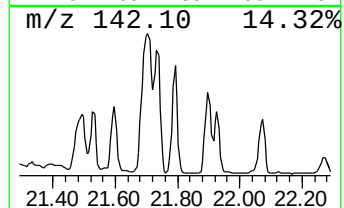
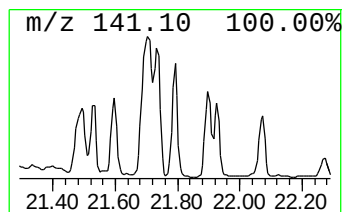
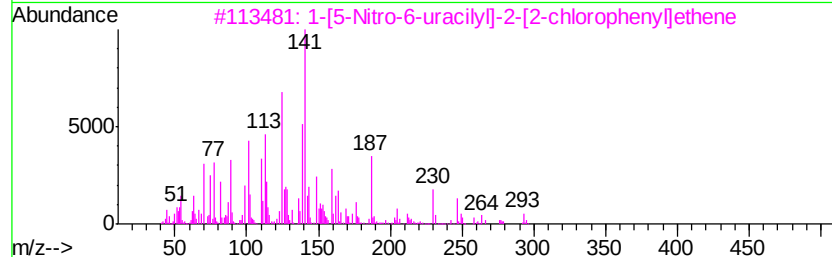
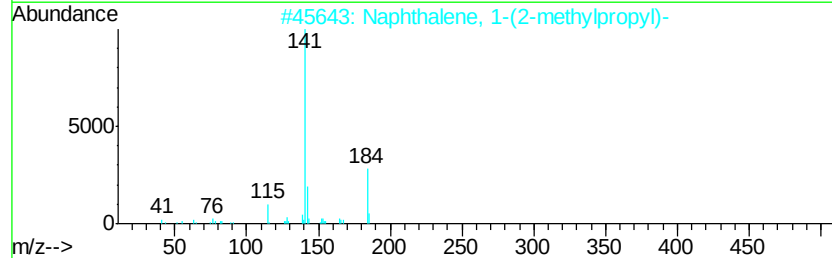
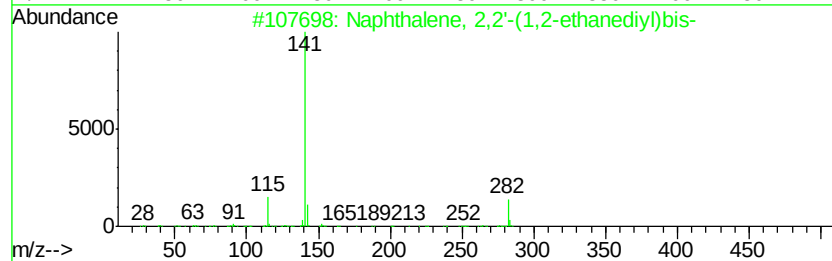
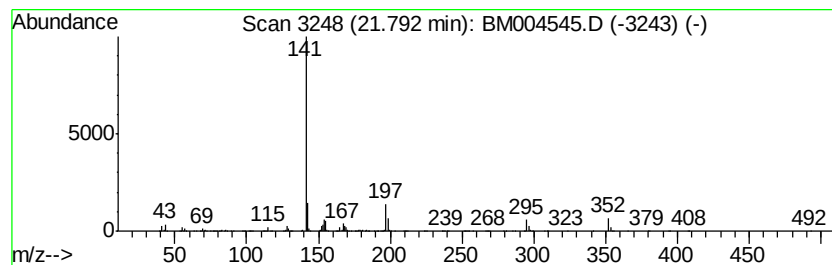
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-04 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.79	23.88 ng/ul	1777880	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,2'-(1,2-ethanediyl...	282	C22H18	021969-45-9	36
2		Naphthalene, 1-(2-methylpropyl)-	184	C14H16	016727-91-6	36
3		1-[5-Nitro-6-uracilyl]-2-[2-chlo...	293	C12H8ClN3O4	296798-53-3	28
4		Naphthalene, 2-pyrrolidino-	197	C14H15N	013672-14-5	10
5		2,4,6(1H,3H,5H)-Pyrimidinetrione...	184	C8H12N2O3	001953-33-9	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

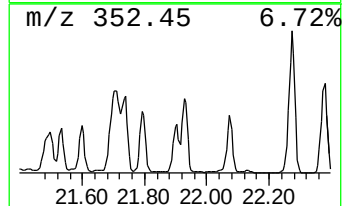
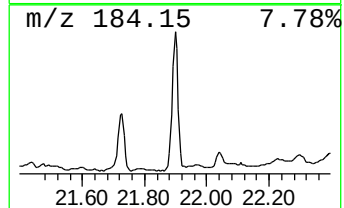
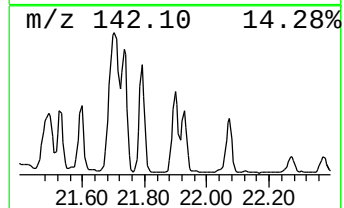
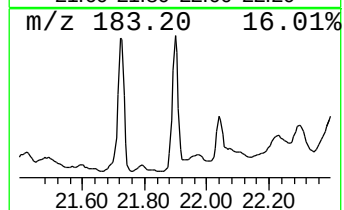
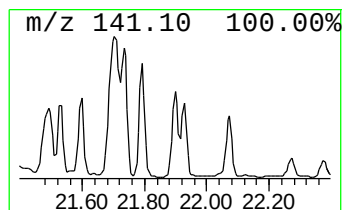
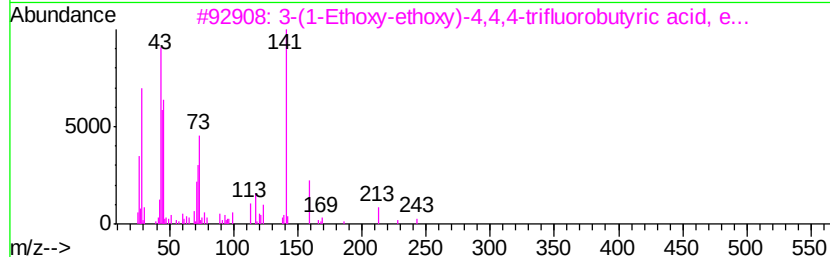
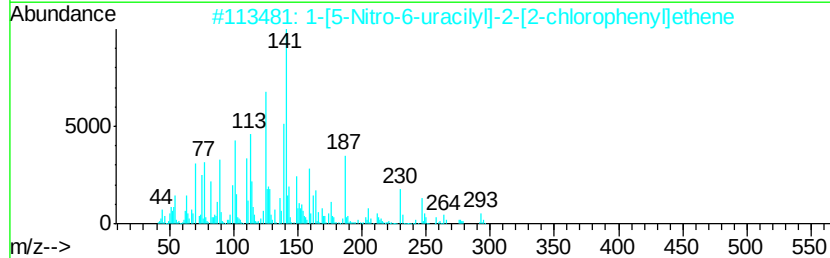
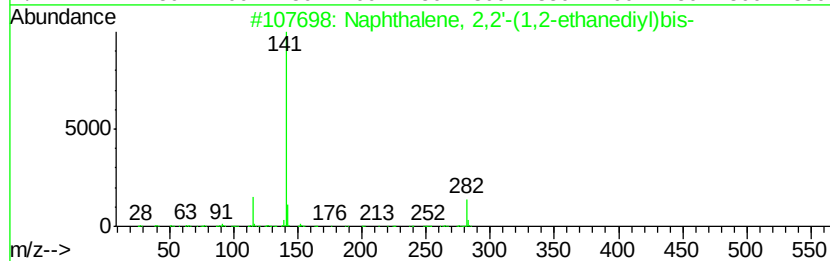
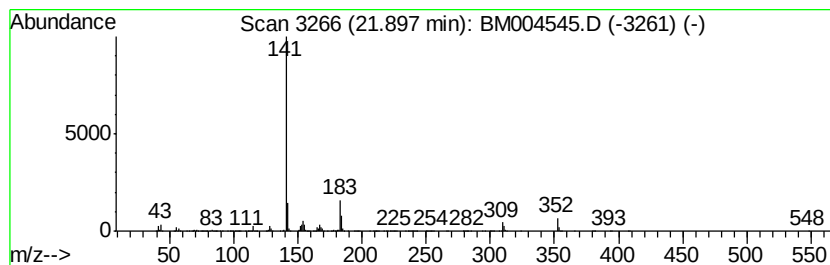
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ClientSampleId :
BLDG06-P003-SS001-01

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Naphthalene, 2,2'-(1,2-etha... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.90	17.68 ng/ul	1315900	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2,2'-(1,2-ethanediy...	282 C22H18	021969-45-9 58
2		1-[5-Nitro-6-uracilyl]-2-[2-chlo...	293 C12H8ClN3O4	296798-53-3 28
3		3-(1-Ethoxy-ethoxy)-4,4,4-triflu...	258 C10H17F3O4	095605-52-0 10
4		Naphthalene, 2-butyl-	184 C14H16	001134-62-9 9
5		Naphthalene, 2-butyl-	184 C14H16	001134-62-9 9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
BLDG06-P003-SS001-01

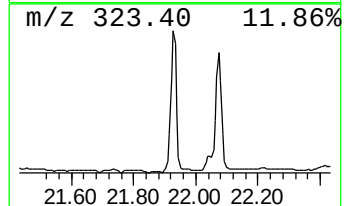
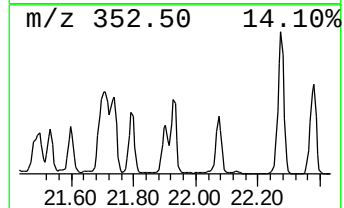
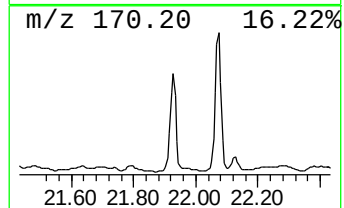
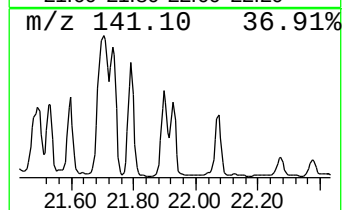
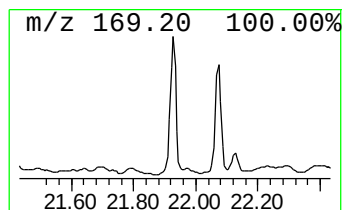
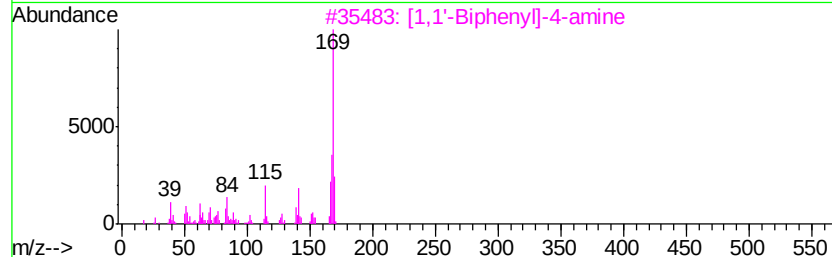
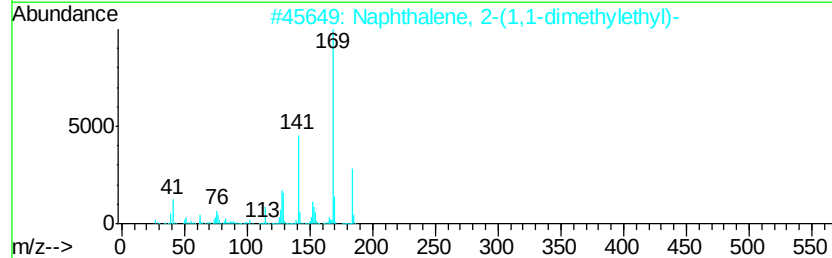
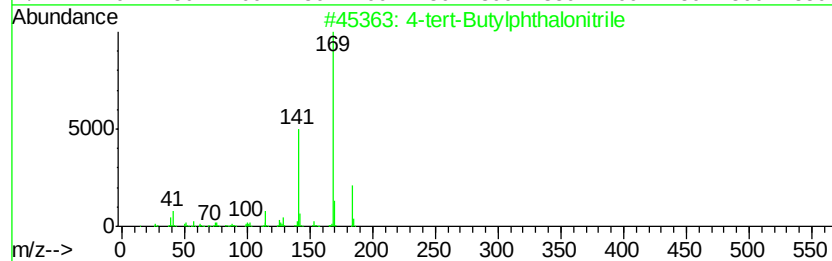
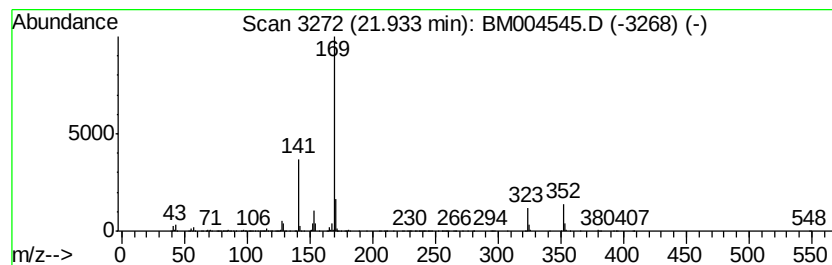
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 4-tert-Butylphthalonitrile Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.93	17.12 ng/ul	1274540	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-tert-Butylphthalonitrile	184	C12H12N2	032703-80-3	59
2		Naphthalene, 2-(1,1-dimethylethyl)-	184	C14H16	002876-35-9	53
3		[1,1'-Biphenyl]-4-amine	169	C12H11N	000092-67-1	47
4		[1,1'-Biphenyl]-3-amine	169	C12H11N	002243-47-2	43
5		4-(4-Methylphenyl)pyridine	169	C12H11N	004423-10-3	43



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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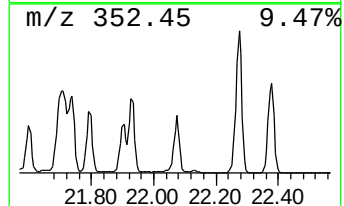
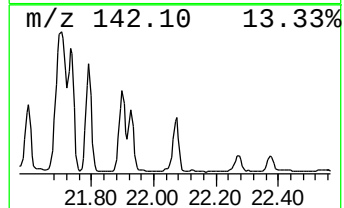
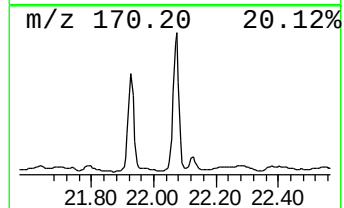
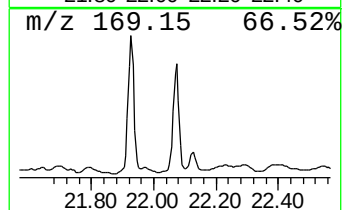
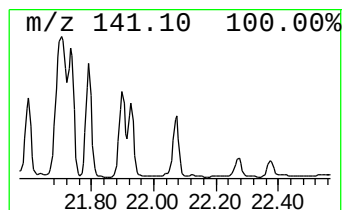
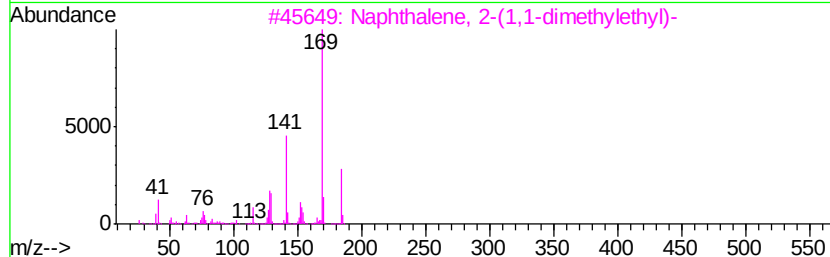
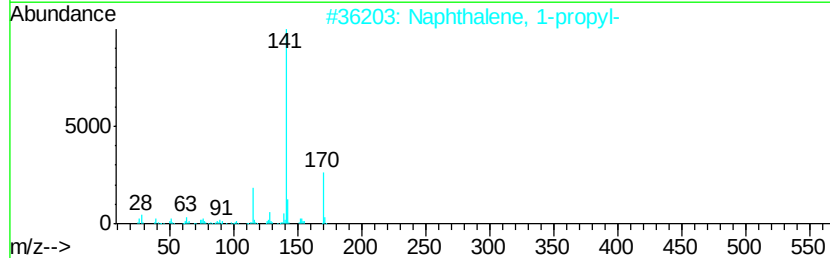
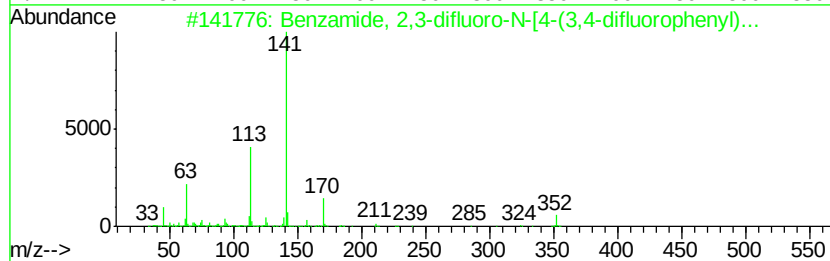
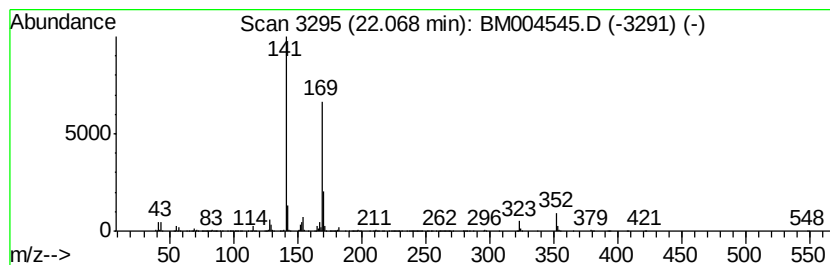
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzamide, 2,3-difluoro-N-[... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.07	16.57 ng/ul	1233620	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzamide, 2,3-difluoro-N-[4-(3,...	352	C16H8F4N2OS	1000277-49-1	50
2		Naphthalene, 1-propyl-	170	C13H14	002765-18-6	38
3		Naphthalene, 2-(1,1-dimethylethyl)-	184	C14H16	002876-35-9	37
4		4-tert-Butylphthalonitrile	184	C12H12N2	032703-80-3	25
5		1-[5-Nitro-6-uracilyl]-2-[2-chlo...	293	C12H8ClN3O4	296798-53-3	17



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
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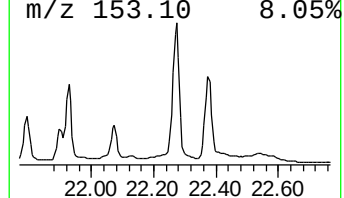
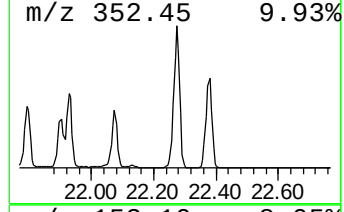
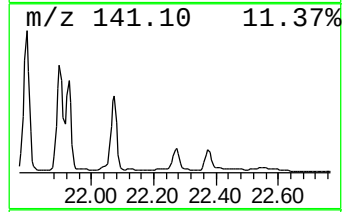
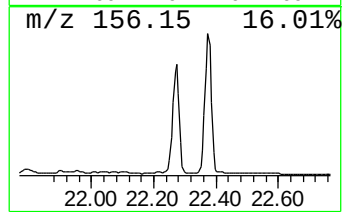
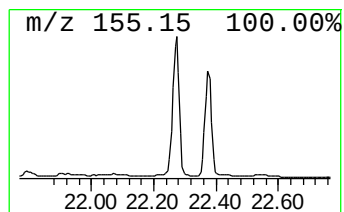
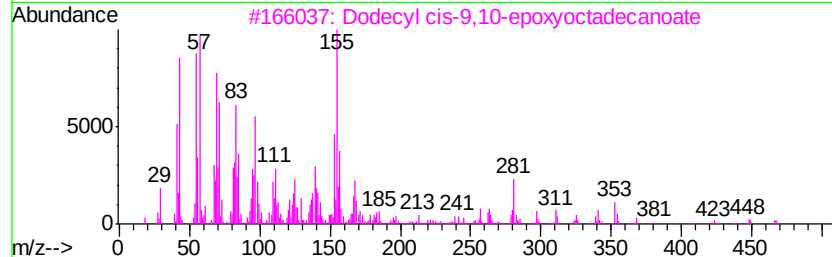
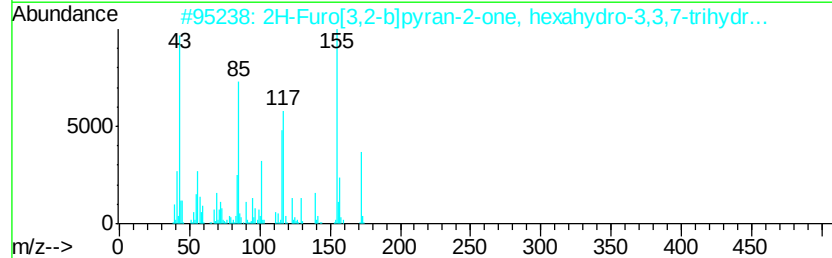
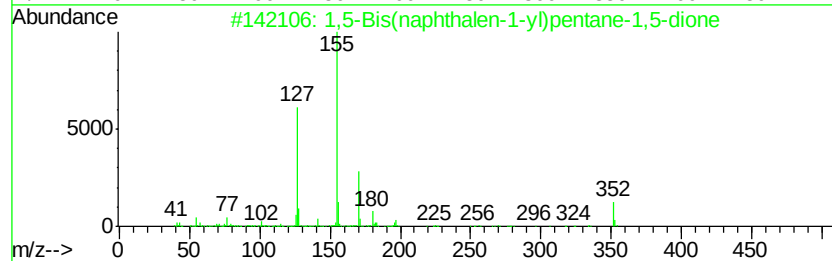
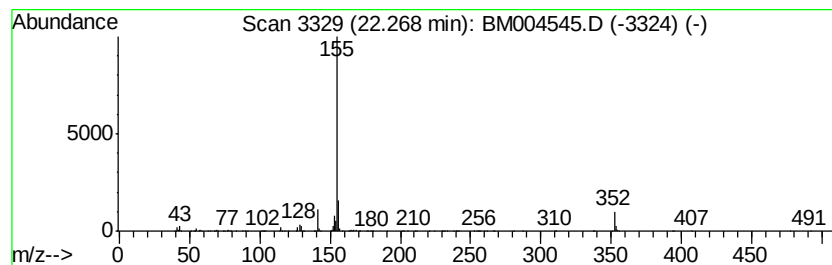
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 1,5-Bis(naphthalen-1-yl)pen... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
22.27	37.29 ng/ul	2776500	Chrysene-d12	21.23		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1,5-Bis(naphthalen-1-yl)pentane-...	352	C25H20O2	1000210-52-9	72
2		2H-Furo[3,2-b]pyran-2-one, hexah...	262	C11H18O7	050392-34-2	33
3		Dodecyl cis-9,10-epoxyoctadecanoate	466	C30H58O3	092332-53-1	16
4		Pyridine, 3-phenyl-	155	C11H9N	001008-88-4	9
5		1,1,1,7,7,7-Hexamethyl-4,4-divin...	284	C14H32Si3	080153-54-4	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
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BLDG06-P003-SS001-01

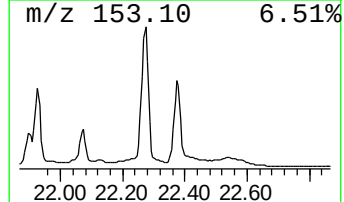
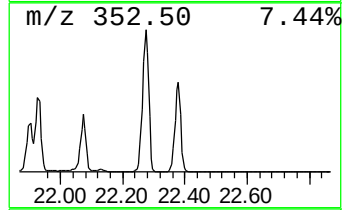
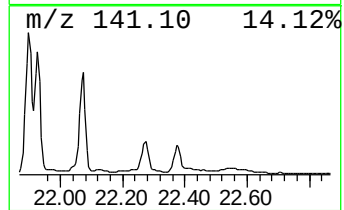
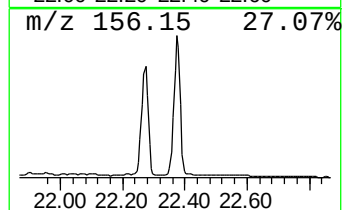
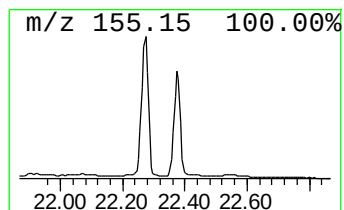
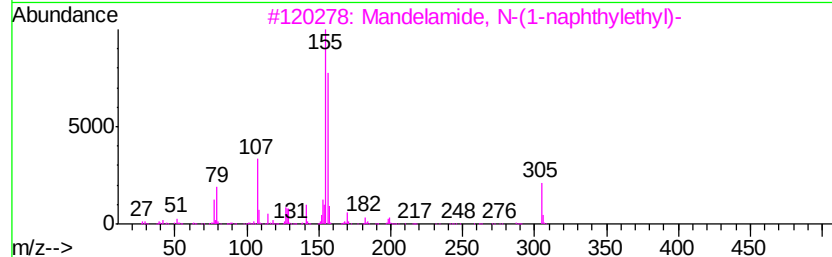
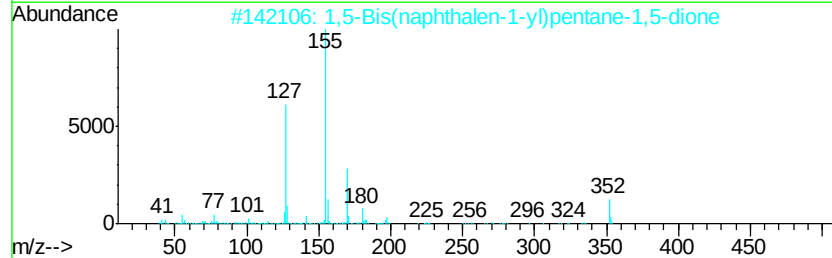
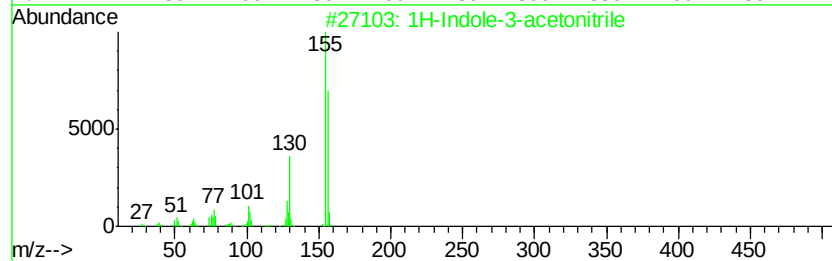
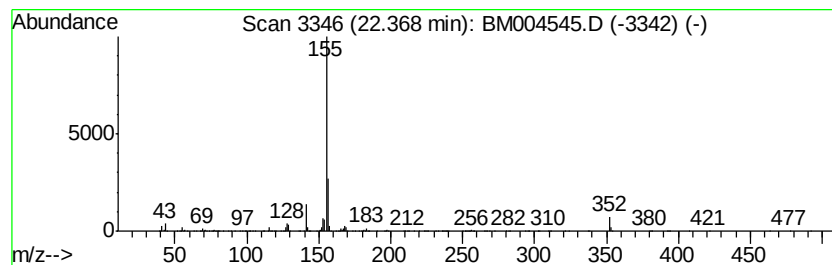
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 unknown-05 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.37	100.63 ng/ul	2101630	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indole-3-acetonitrile	156	C10H8N2	000771-51-7	10
2		1,5-Bis(naphthalen-1-yl)pentane-...	352	C25H20O2	1000210-52-9	10
3		Mandelamide, N-(1-naphthylethyl)-	305	C20H19NO2	344875-77-0	10
4		1H-Indole-3-acetonitrile	156	C10H8N2	000771-51-7	9
5		6-Amino-1,3-dimethyluracil	155	C6H9N3O2	006642-31-5	9



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030316\
Data File : BM004545.D
Acq On : 03 Mar 2016 23:13
Operator : SJ/UM
Sample : H1616-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BLDG06-P003-SS001-01

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM030316.M
Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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(DEL) Alkane: Str...	21.53	10.6	ng/ul	786053	5	21.23	1488940	20.0
unknown-02	21.60	15.1	ng/ul	1120730	5	21.23	1488940	20.0
Naphthalene, 1-(2...	21.70	47.4	ng/ul	3526240	5	21.23	1488940	20.0
unknown-03	21.73	23.3	ng/ul	1734030	5	21.23	1488940	20.0
unknown-04	21.79	23.9	ng/ul	1777880	5	21.23	1488940	20.0
Naphthalene, 2,2'...	21.90	17.7	ng/ul	1315900	5	21.23	1488940	20.0
4-tert-Butylphtha...	21.93	17.1	ng/ul	1274540	5	21.23	1488940	20.0
Benzamide, 2,3-di...	22.07	16.6	ng/ul	1233620	5	21.23	1488940	20.0
1,5-Bis(naphthale...	22.27	37.3	ng/ul	2776500	5	21.23	1488940	20.0
unknown-05	22.37	100.6	ng/ul	2101630	6	23.44	417687	20.0