

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110116\
 Data File : BG024613.D
 Acq On : 2 Nov 2016 5:05
 Operator : UM/SJ
 Sample : H5489-20
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-77-0.5-1.0

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % 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\8270-BG102516.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.131	44	50	67	rVB	5384575	9793565	100.00%	17.511%
2	5.328	419	424	433	rBV	426332	689692	7.04%	1.233%
3	5.798	496	504	513	rBV	1884088	3212253	32.80%	5.744%
4	7.402	769	777	786	rBV	2051647	3631234	37.08%	6.493%
5	7.796	836	844	852	rVB	1865401	3357972	34.29%	6.004%
6	8.266	918	924	935	rVB	426150	775828	7.92%	1.387%
7	8.595	971	980	991	rVB	1344432	2405606	24.56%	4.301%
8	9.441	1115	1124	1133	rBV	1253881	2284980	23.33%	4.086%
9	11.092	1397	1405	1415	rVB	552367	1026267	10.48%	1.835%
10	13.507	1807	1816	1824	rBV	2980286	4739176	48.39%	8.474%
11	14.318	1948	1954	1963	rBV	742363	1135927	11.60%	2.031%
12	14.882	2043	2050	2057	rBV	955415	1518996	15.51%	2.716%
13	16.363	2295	2302	2312	rBV	2767544	4358329	44.50%	7.793%
14	16.539	2328	2332	2339	rBV2	76453	135803	1.39%	0.243%
15	17.626	2510	2517	2527	rVB	1231648	1933611	19.74%	3.457%
16	17.961	2568	2574	2581	rBV4	118520	216045	2.21%	0.386%
17	18.466	2653	2660	2666	rBV3	110715	173382	1.77%	0.310%
18	20.205	2948	2956	2962	rBV	6127927	8695775	88.79%	15.548%
19	21.504	3172	3177	3185	rBV7	103746	220177	2.25%	0.394%
20	21.921	3241	3248	3255	rVB	1509090	2729519	27.87%	4.880%
21	23.648	3535	3542	3546	rBV3	109695	200162	2.04%	0.358%
22	25.346	3819	3831	3845	rVB	872732	2693765	27.51%	4.816%

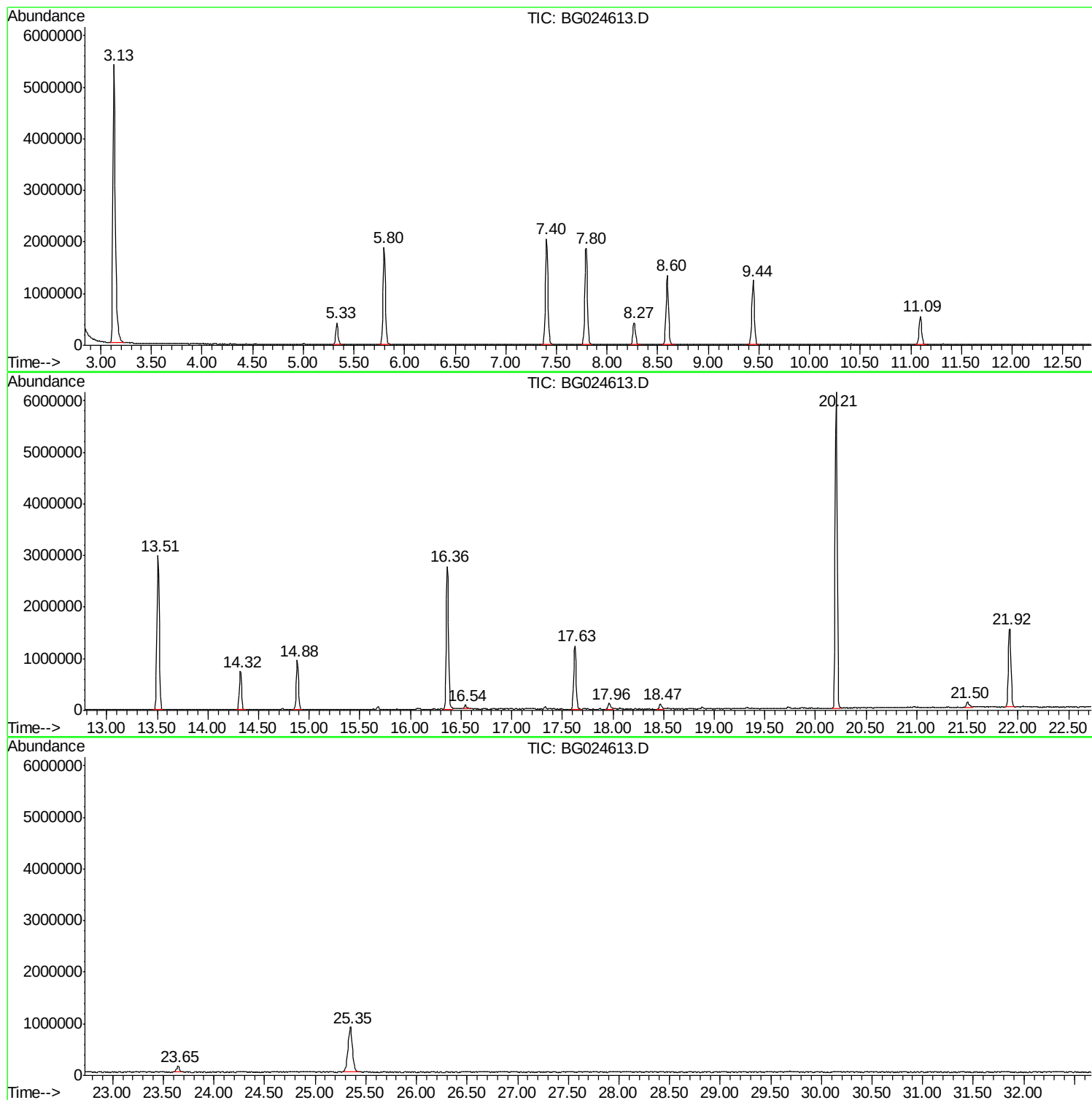
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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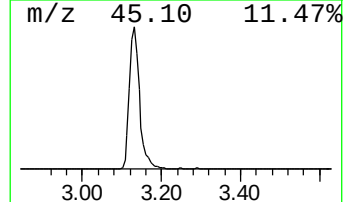
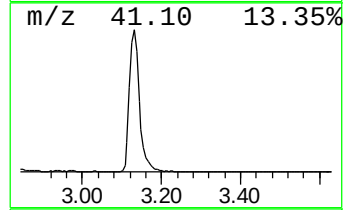
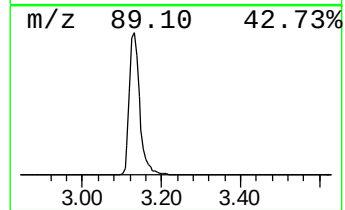
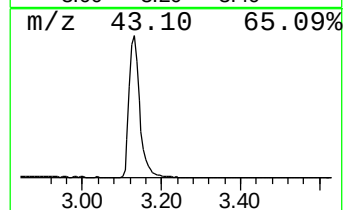
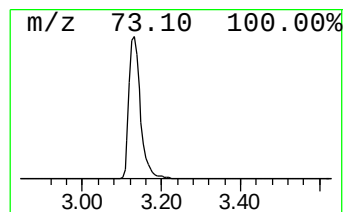
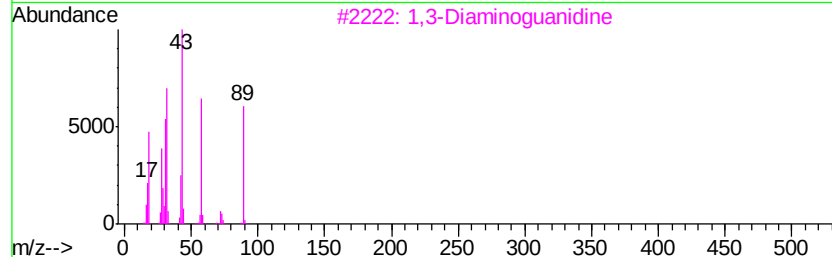
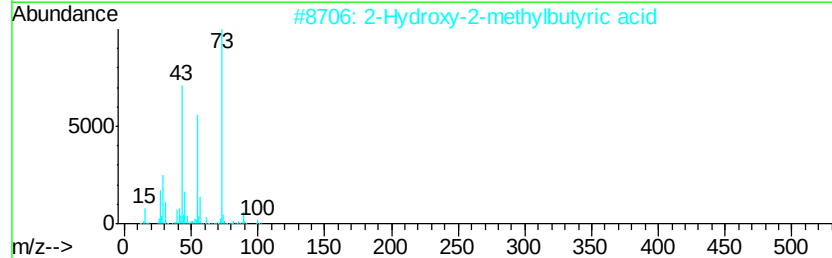
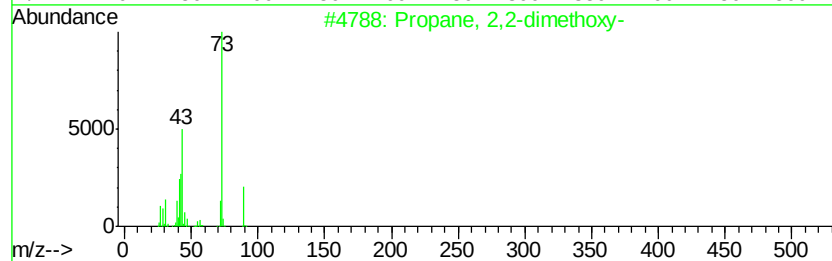
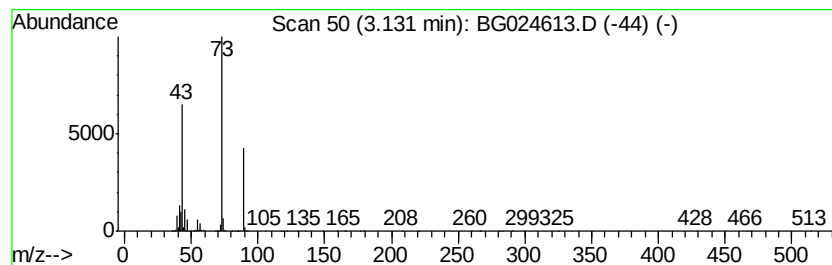
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown3.13 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.13	252.47 ng	9793570	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	40
2		2-Hydroxy-2-methylbutyric acid	118	C5H10O3	003739-30-8	17
3		1,3-Diaminoguanidine	89	CH7N5	004364-78-7	9
4		N-Ethylformamide	73	C3H7NO	000627-45-2	9
5		1-Hexen-4-ol, 1-chloro-3,5-dimet...	162	C8H15ClO	100244-09-5	9



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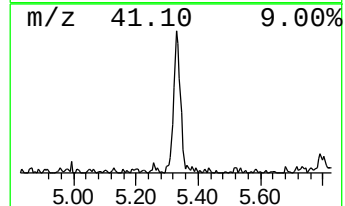
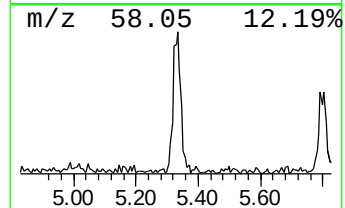
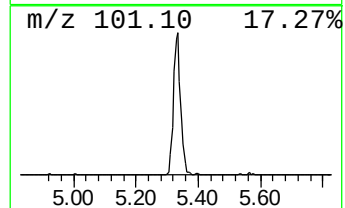
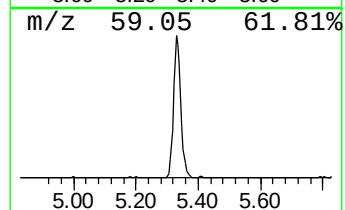
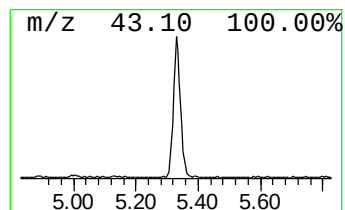
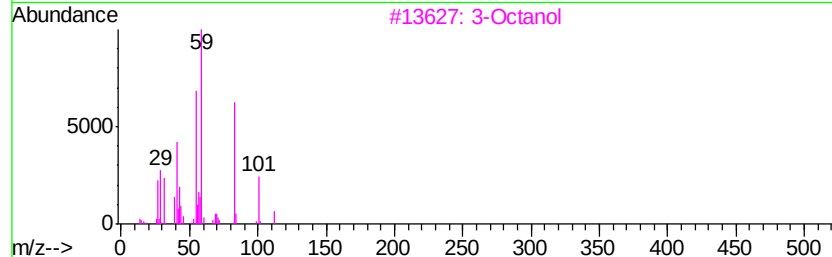
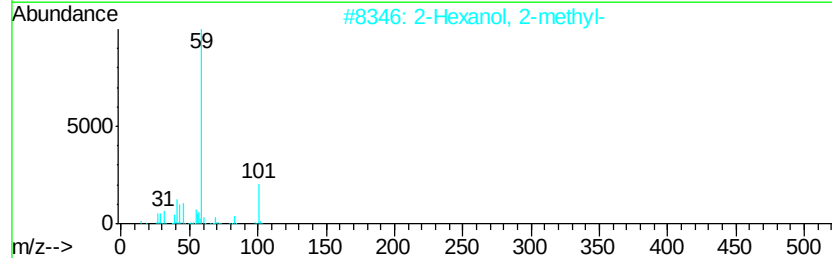
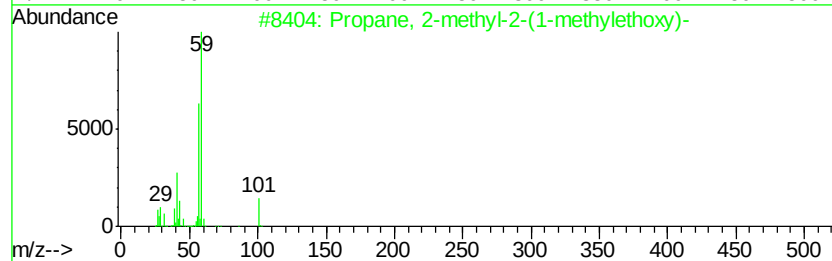
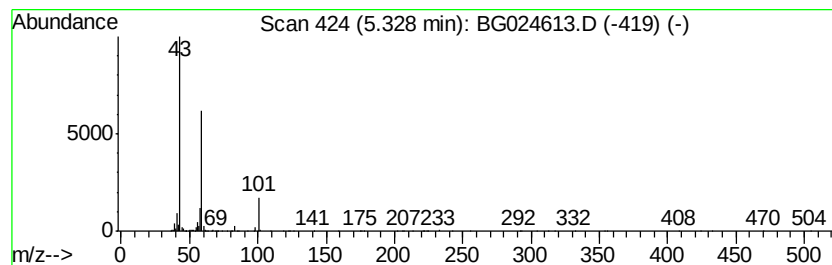
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 unknown5.33 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.33	17.78 ng	689692	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	38
2		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	37
3		3-Octanol	130	C8H18O	000589-98-0	33
4		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	32
5		Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	17



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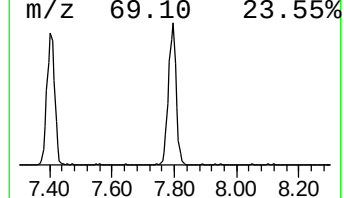
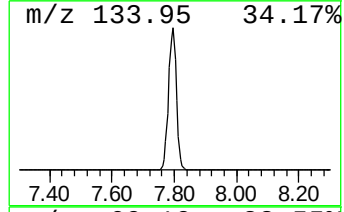
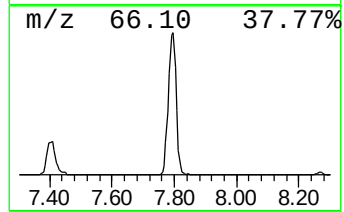
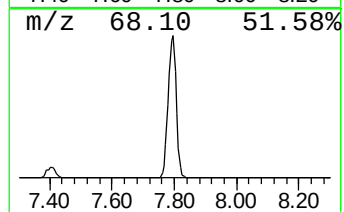
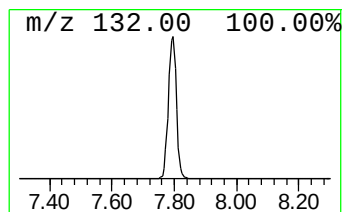
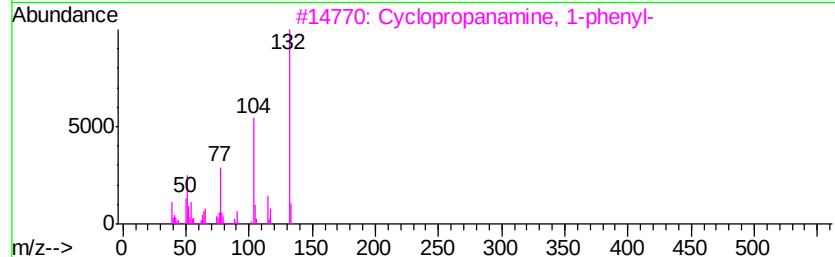
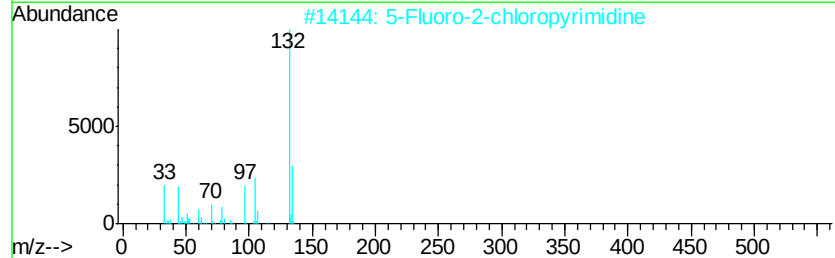
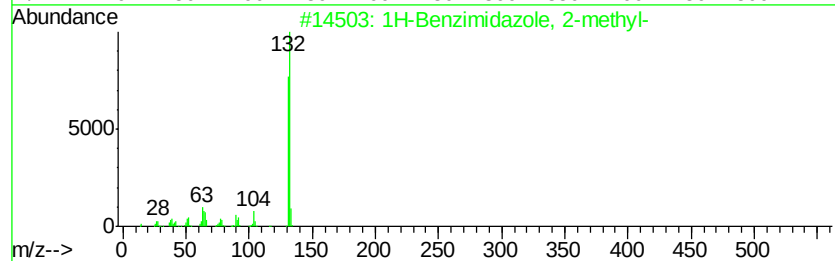
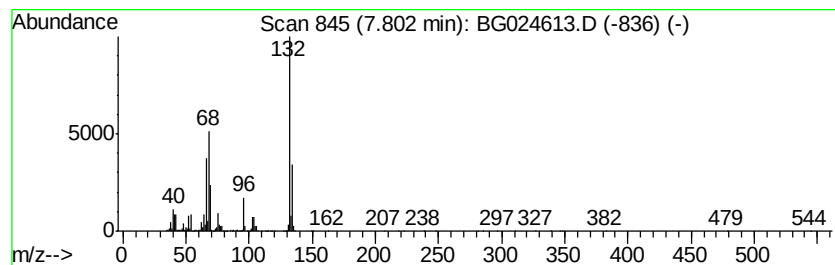
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Peak Number 3 unknown7.80 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.80	86.56 ng	3357970	1,4-Dichlorobenzene-d4	8.27

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	10
3		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10



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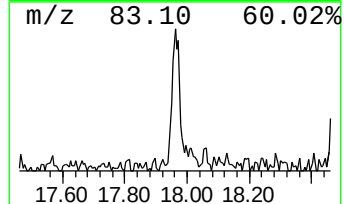
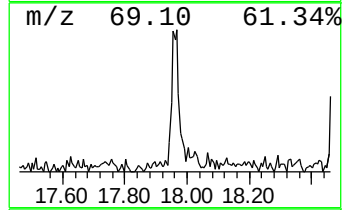
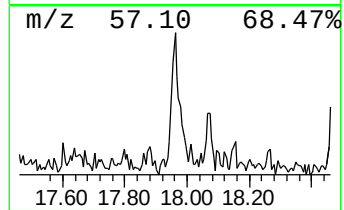
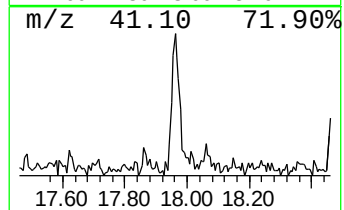
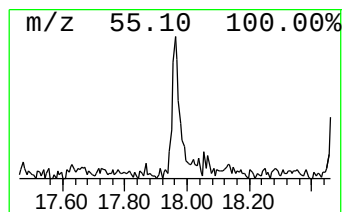
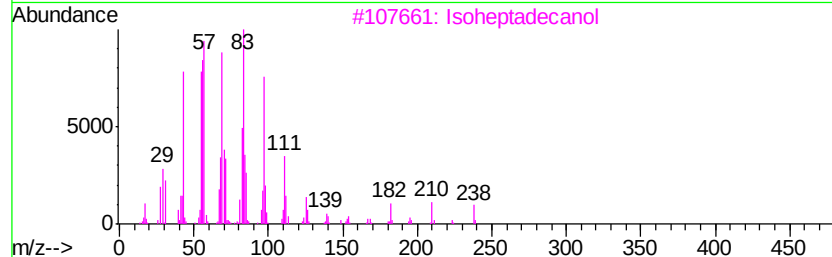
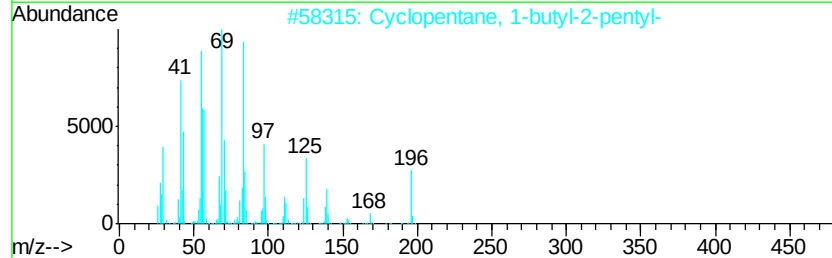
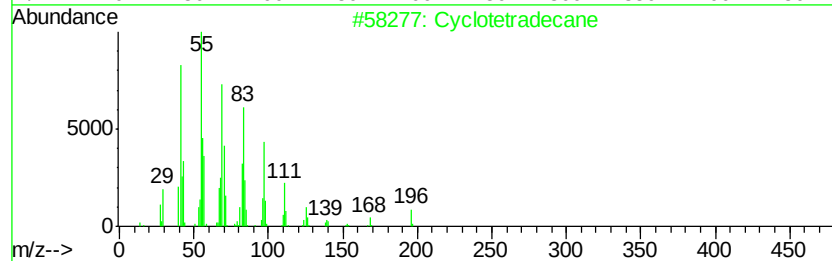
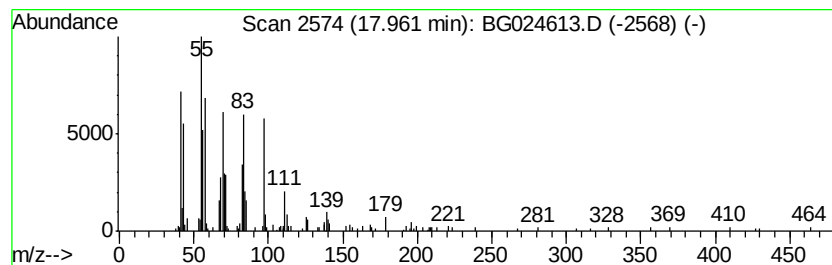
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Peak Number 5 Cyclotetradecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.96	2.23 ng	216045	Phenanthrene-d10	17.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	86
2	Cyclopentane, 1-butyl-2-pentyl-	196	C14H28	061142-52-7	76
3	Isoheptadecanol	256	C17H36O	057289-07-3	72
4	1-Octadecanol	270	C18H38O	000112-92-5	72
5	9-Octadecene, (E)-	252	C18H36	007206-25-9	64



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
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unknown5.33	5.33	17.8	ng	689692	1	8.27	775828	20.0
unknown7.80	7.80	86.6	ng	3357970	1	8.27	775828	20.0
Cyclotetradecane	17.96	2.2	ng	216045	4	17.63	1933610	20.0