

Data Path : Z:\HPCHEM1\BNA G\DATA\BG092816\
 Data File : BG024154.D
 Acq On : 28 Sep 2016 20:47
 Operator : UM/SJ
 Sample : H4992-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 092716-DRDC-F-D-M

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-BG092316.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	5.106	383	386	392	rVB2	45063	66046	1.27%	0.341%
2	5.588	464	468	477	rBV	220941	387894	7.48%	2.005%
3	7.215	740	745	759	rBV2	164923	360965	6.96%	1.866%
4	7.567	799	805	814	rBV2	518610	917184	17.70%	4.741%
5	8.037	879	885	894	rVB	186610	321328	6.20%	1.661%
6	8.360	933	940	949	rBV	675592	1208334	23.31%	6.246%
7	9.218	1079	1086	1100	rBV	631827	1232024	23.77%	6.368%
8	10.869	1360	1367	1381	rBV	236037	474519	9.16%	2.453%
9	13.324	1777	1785	1794	rBV	1837098	3015470	58.18%	15.587%
10	14.164	1922	1928	1942	rBV2	90294	166729	3.22%	0.862%
11	14.711	2014	2021	2032	rBV	463528	782218	15.09%	4.043%
12	16.215	2271	2277	2291	rBV	1038778	1754254	33.85%	9.068%
13	17.483	2486	2493	2506	rBV2	637631	1047709	20.22%	5.416%
14	18.353	2635	2641	2646	rBV6	37243	67385	1.30%	0.348%
15	19.622	2850	2857	2866	rBV6	35715	78382	1.51%	0.405%
16	20.086	2930	2936	2947	rBV	3972080	5182654	100.00%	26.789%
17	21.783	3219	3225	3238	rVB	663941	1219431	23.53%	6.303%
18	25.120	3783	3793	3807	rBV2	348930	1063340	20.52%	5.496%

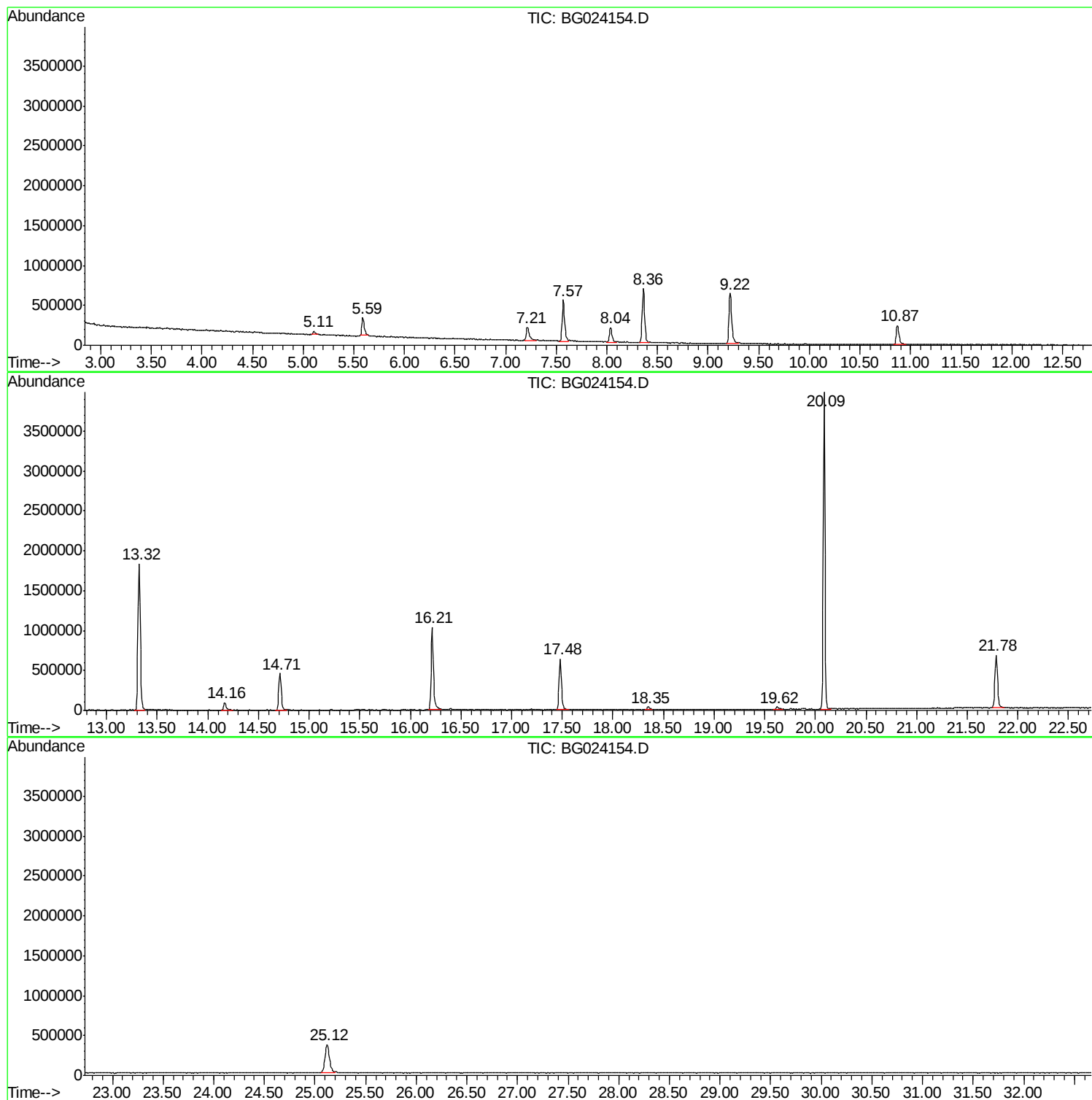
Sum of corrected areas: 19345866

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG092316.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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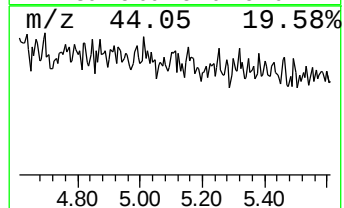
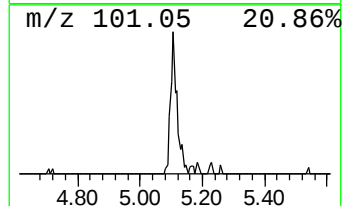
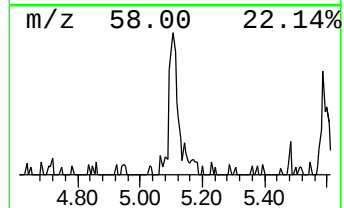
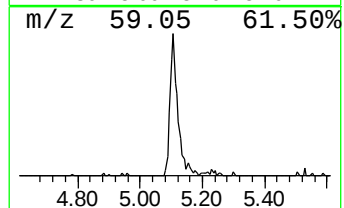
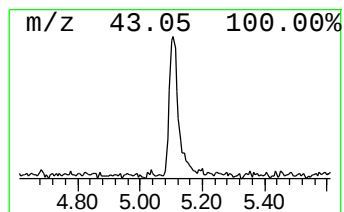
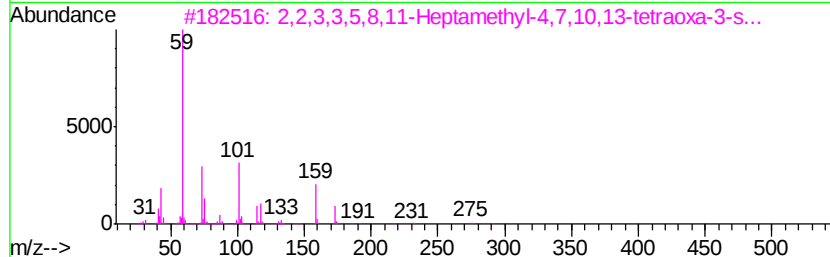
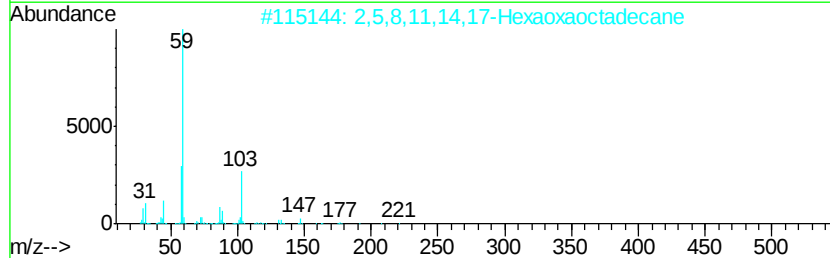
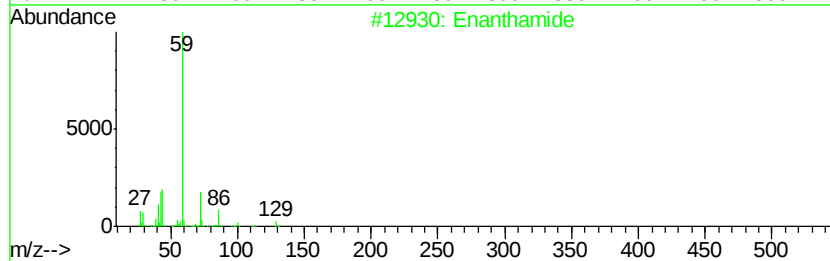
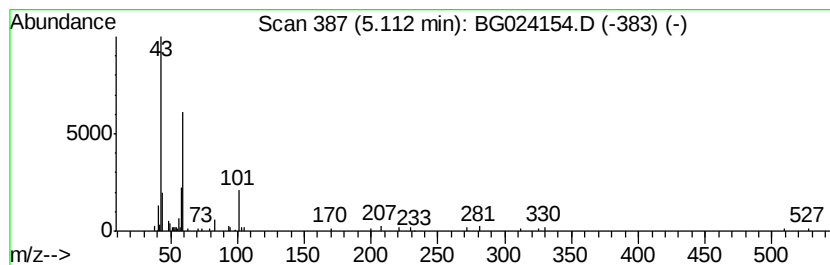
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown5.11 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	4.11 ng	66046	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Enanthamide	129	C7H15NO	000628-62-6	37
2			2,5,8,11,14,17-Hexaoxaoctadecane	266	C12H26O6	001191-87-3	36
3			2,2,3,3,5,8,11-Heptamethyl-4,7,1...	348	C18H40O4Si	1000367-07-2	36
4			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	33
5			2-Butanol, 3-methoxy-	104	C5H12O2	053778-72-6	28



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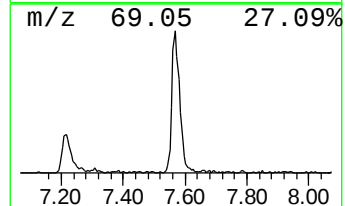
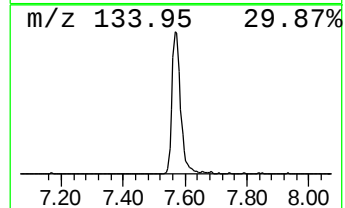
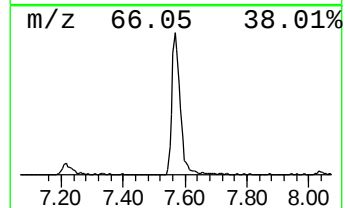
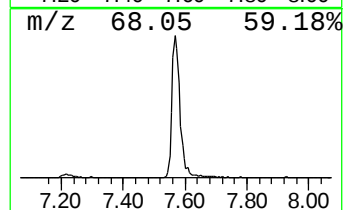
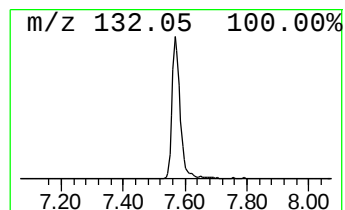
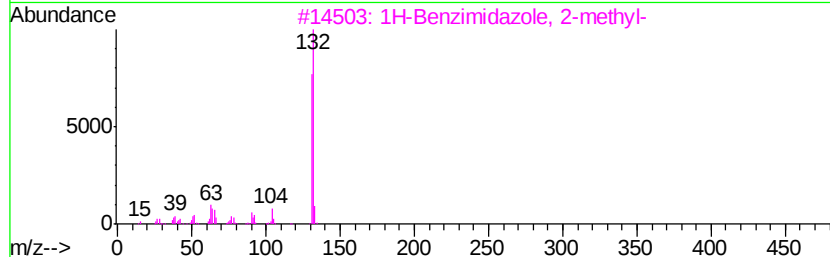
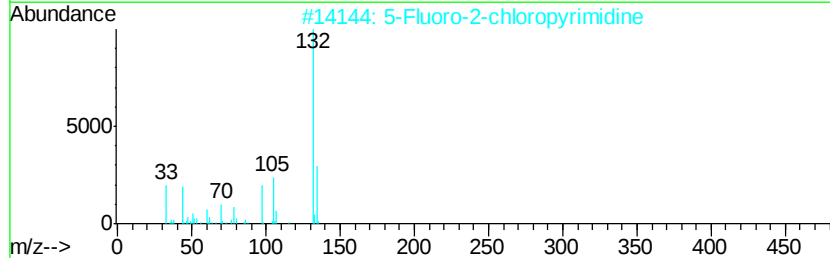
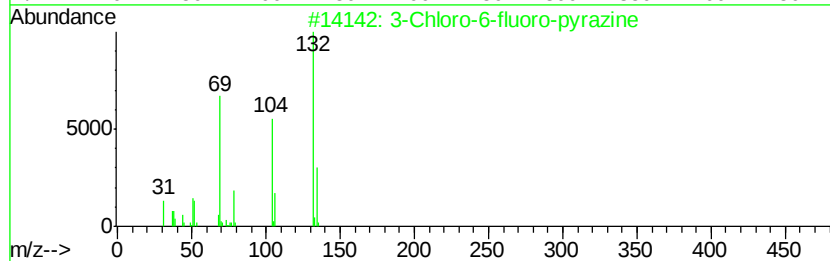
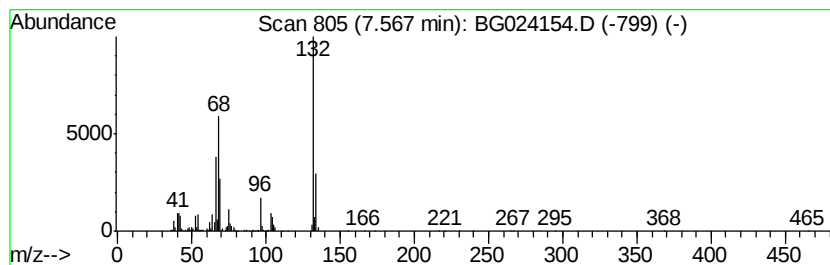
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Peak Number 2 unknown7.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	57.09 ng	917184	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	30
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	12
4		Pyrazolo(2,3-a)pyridine, 7-methyl-	132	C8H8N2	034760-58-2	12
5		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown5.11	5.11	4.1	ng	66046	1	8.04	321328	20.0
unknown7.57	7.57	57.1	ng	917184	1	8.04	321328	20.0