

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041816\
 Data File : BF086287.D
 Acq On : 18 Apr 2016 17:42
 Operator : UM/SJ
 Sample : H2413-06 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2-F-6(0-2)

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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_F\METHODS\8270-BF041516.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	5.059	259	260	264	rBV	116364	135417	5.89%	0.599%
2	5.470	294	296	299	rBV	1718072	1604926	69.86%	7.094%
3	6.499	384	386	389	rBV	1712473	1553155	67.61%	6.865%
4	6.647	396	399	401	rBV	2038422	1760499	76.63%	7.781%
5	6.887	417	420	422	rBV	2001723	1817183	79.10%	8.032%
6	7.036	431	433	436	rBV	1159643	1242080	54.07%	5.490%
7	7.447	466	469	471	rBV	1010014	1016513	44.25%	4.493%
8	8.167	530	532	535	rBV	2156126	2252443	98.05%	9.956%
9	9.242	624	626	629	rVB	1205121	1446182	62.95%	6.392%
10	9.333	629	634	635	rBV2	16421	24152	1.05%	0.107%
11	9.642	659	661	663	rBV	187079	157767	6.87%	0.697%
12	9.859	679	680	683	rVB	34664	30541	1.33%	0.135%
13	9.928	683	686	688	rBV	2513577	2297330	100.00%	10.154%
14	10.362	722	724	726	rVB	69752	55084	2.40%	0.243%
15	10.579	740	743	746	rBV	23288	41563	1.81%	0.184%
16	10.705	752	754	757	rBV	844166	769131	33.48%	3.400%
17	10.831	764	765	770	rVB	71517	92847	4.04%	0.410%
18	11.276	802	804	806	rBV	29712	39017	1.70%	0.172%
19	11.402	813	815	818	rBV2	1610672	1840986	80.14%	8.137%
20	11.482	819	822	824	rVB	20847	30436	1.32%	0.135%
21	11.928	858	861	863	rBV2	22018	42056	1.83%	0.186%
22	12.019	867	869	873	rVB	22484	36122	1.57%	0.160%
23	12.202	882	885	890	rBV6	7841	25764	1.12%	0.114%
24	12.614	919	921	923	rVB	125532	102740	4.47%	0.454%
25	12.842	937	941	942	rBV	113365	119379	5.20%	0.528%
26	12.991	952	954	956	rVB	915199	823551	35.85%	3.640%
27	13.894	1030	1033	1037	rBV4	32932	67731	2.95%	0.299%
28	14.042	1043	1046	1048	rBV	1249859	1607744	69.98%	7.106%
29	14.522	1086	1088	1089	rBV	27969	35418	1.54%	0.157%
30	14.888	1118	1120	1123	rBV3	36125	73758	3.21%	0.326%
31	15.060	1133	1135	1139	rVB	48872	108265	4.71%	0.479%
32	15.345	1158	1160	1163	rBV	45884	60490	2.63%	0.267%
33	15.471	1168	1171	1174	rBV	1056738	1314424	57.22%	5.810%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041516.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

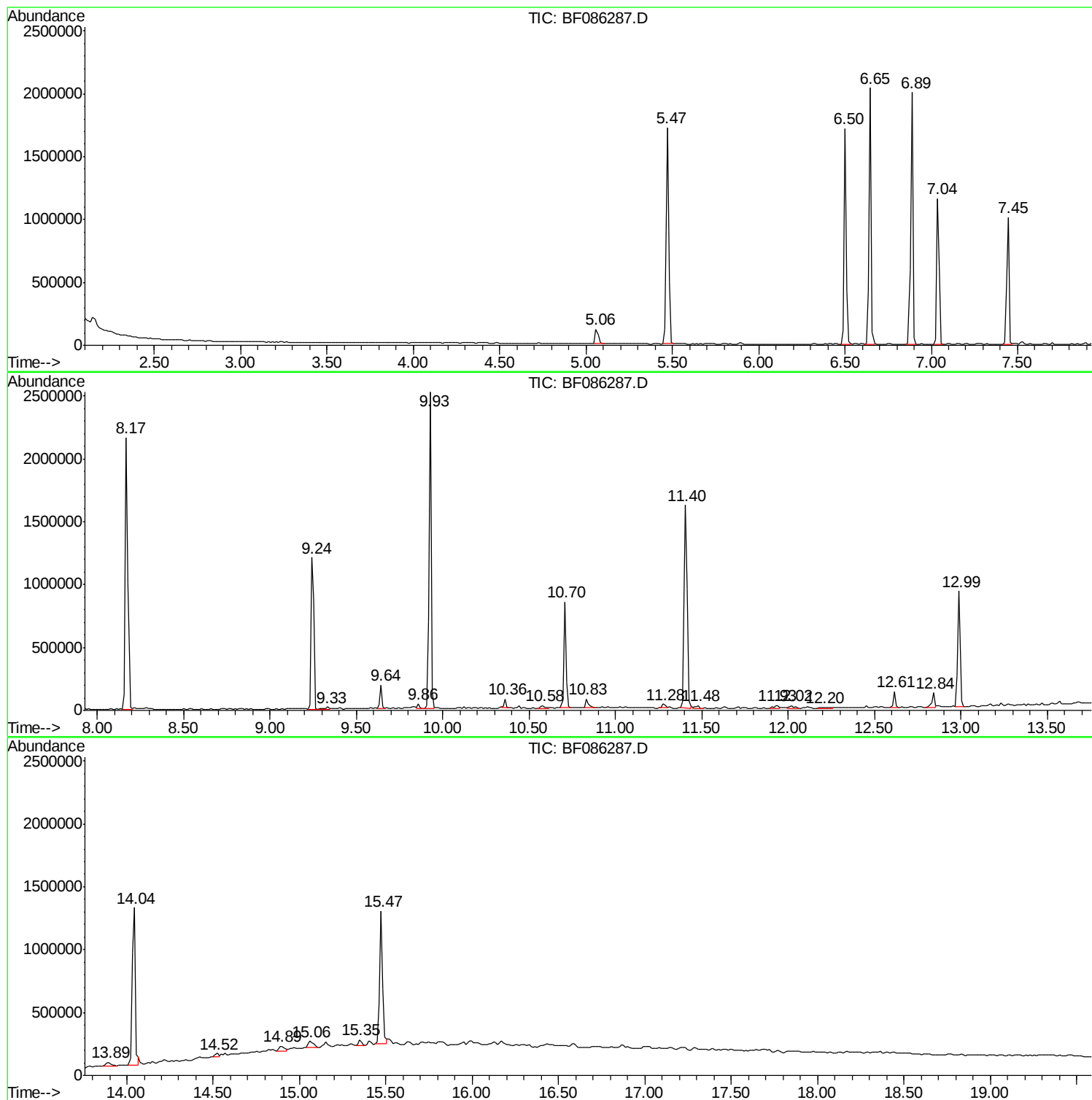
Sum of corrected areas: 22624694

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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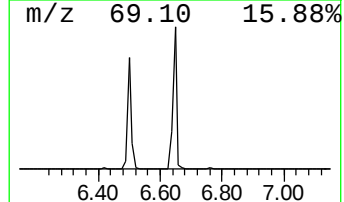
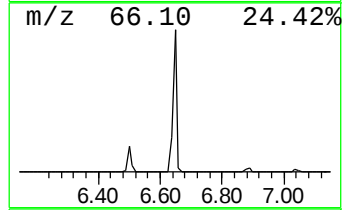
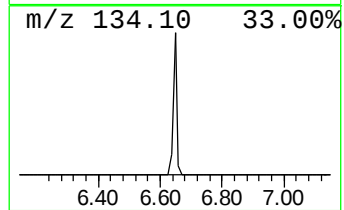
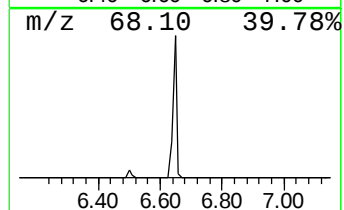
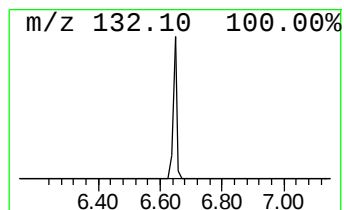
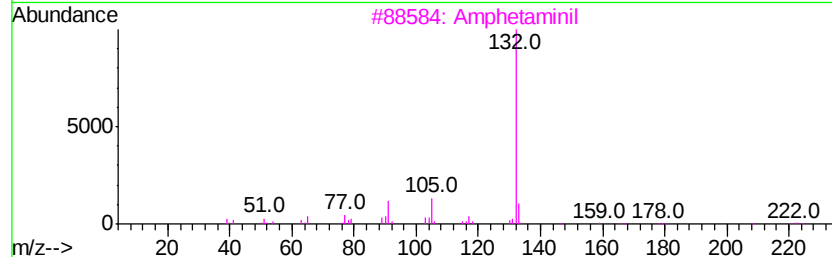
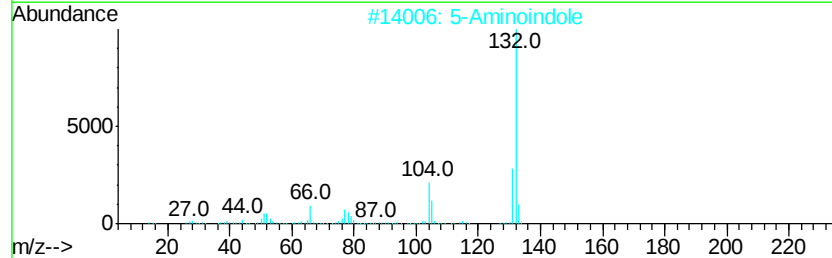
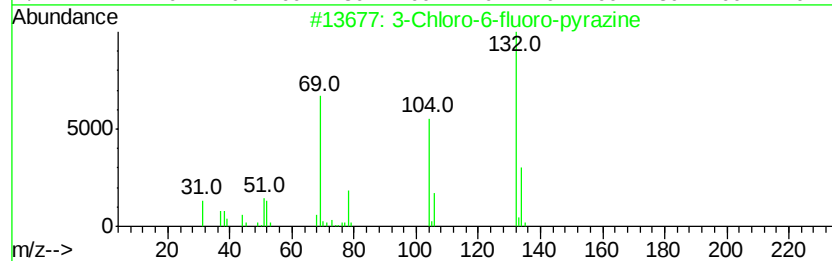
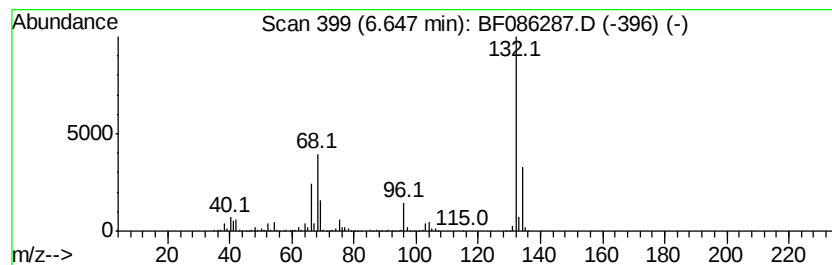
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF041516.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.65 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.65	19.38 ng	1760500	1,4-Dichlorobenzene-d4	6.89

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Amphetaminil	250	C17H18N2	017590-01-1	10
4			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5			Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.65	6.65	19.4	ng	1760500	1	6.89	1817180	20.0