

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF110217\
 Data File : BF100088.D
 Acq On : 2 Nov 2017 16:46
 Operator : SJ/JU
 Sample : I6236-04
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FM782

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-BF102317.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	4.981	490	492	511	rBV	45071	73404	1.36%	0.334%
2	5.393	559	562	581	rBV	468006	683557	12.64%	3.114%
3	6.428	735	738	754	rBV	247441	461289	8.53%	2.101%
4	6.540	754	757	762	rVV	1355367	1476951	27.31%	6.728%
5	6.575	762	763	766	rVV2	137481	126995	2.35%	0.578%
6	6.604	766	768	781	rVV2	121414	183345	3.39%	0.835%
7	6.698	781	784	793	rVV	185117	215353	3.98%	0.981%
8	6.769	793	796	803	rVV	546902	479069	8.86%	2.182%
9	6.922	818	822	833	rBV	1696798	1598834	29.56%	7.283%
10	7.340	890	893	910	rVB	1095143	1173040	21.69%	5.343%
11	7.898	979	988	991	rBV	3109711	5407901	100.00%	24.634%
12	8.051	1009	1014	1024	rVB	784524	730164	13.50%	3.326%
13	8.269	1048	1051	1053	rBV	468067	360820	6.67%	1.644%
14	8.292	1053	1055	1059	rVV	519179	433405	8.01%	1.974%
15	8.416	1073	1076	1084	rVB	239398	246328	4.55%	1.122%
16	8.498	1087	1090	1091	rBV	98511	87623	1.62%	0.399%
17	8.516	1091	1093	1097	rVV2	161175	165500	3.06%	0.754%
18	9.128	1193	1197	1200	rBV	2613388	2326677	43.02%	10.598%
19	9.528	1262	1265	1268	rVV	128584	113018	2.09%	0.515%
20	9.810	1309	1313	1316	rBV	771158	714939	13.22%	3.257%
21	10.604	1444	1448	1455	rBV	1191745	1124210	20.79%	5.121%
22	11.304	1563	1567	1575	rBV	722268	682413	12.62%	3.109%
23	11.822	1651	1655	1658	rBV	188527	177565	3.28%	0.809%
24	12.527	1769	1775	1779	rBV5	27781	58356	1.08%	0.266%
25	12.604	1785	1788	1796	rBV	182183	187623	3.47%	0.855%
26	12.774	1814	1817	1825	rVB	133640	135642	2.51%	0.618%
27	12.886	1832	1836	1840	rBV	1712515	1552659	28.71%	7.073%
28	13.786	1985	1989	1994	rBV2	44032	54940	1.02%	0.250%
29	13.980	2018	2022	2027	rBV	490801	460891	8.52%	2.099%
30	15.504	2277	2281	2291	rVB	352759	460457	8.51%	2.097%

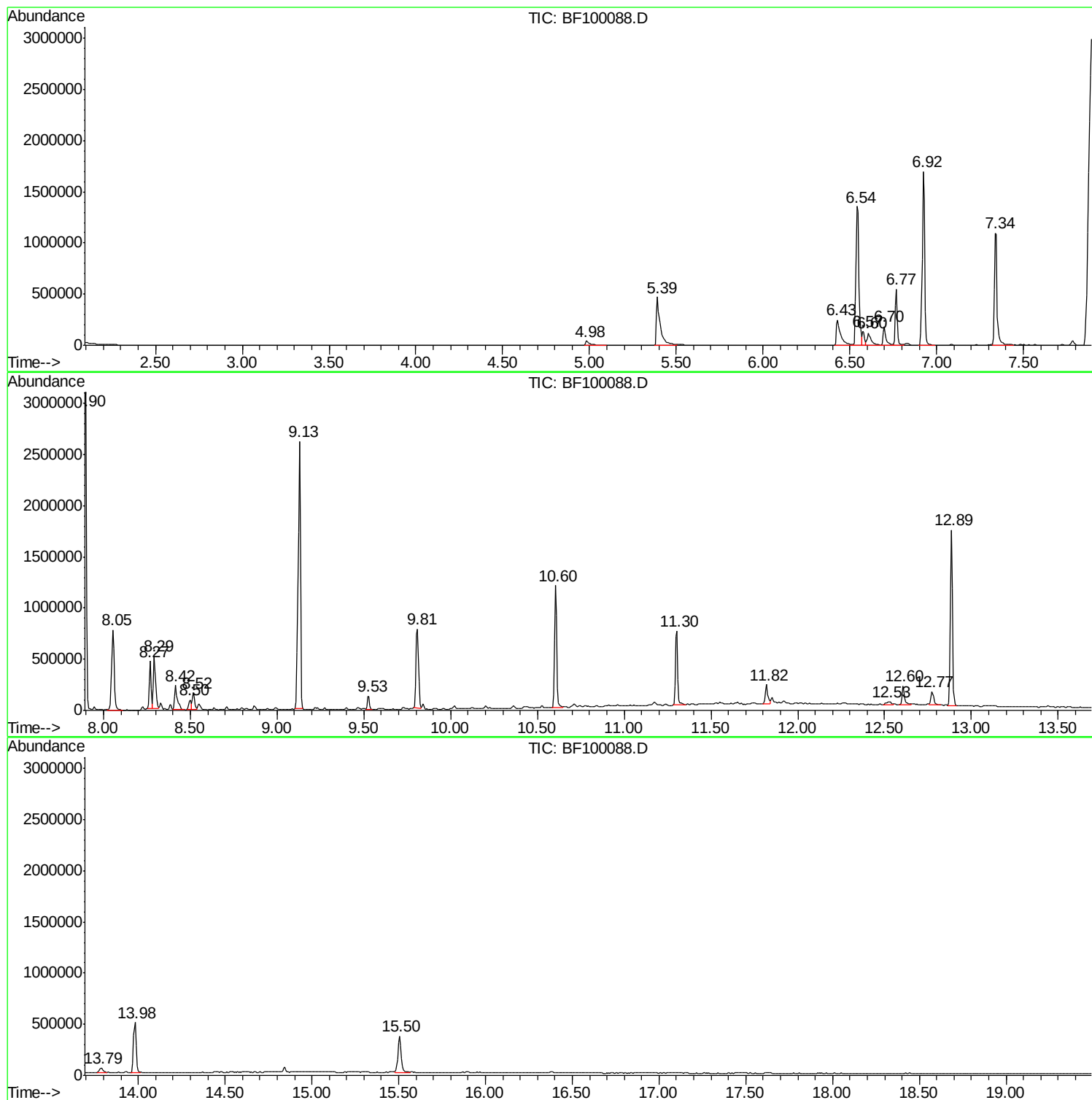
Sum of corrected areas: 21952968

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ClientSampled :
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.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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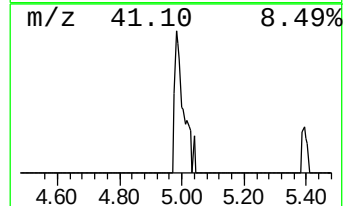
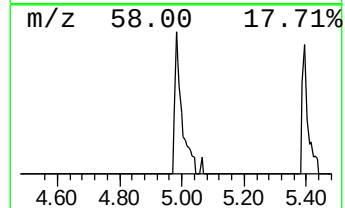
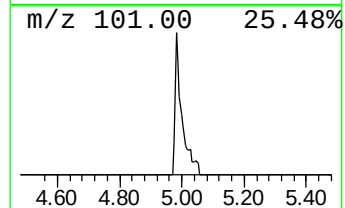
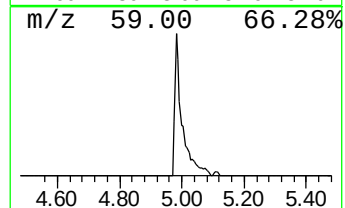
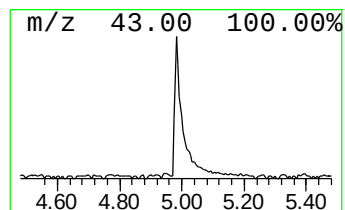
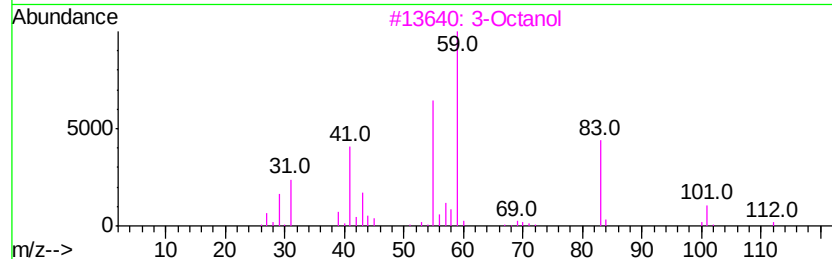
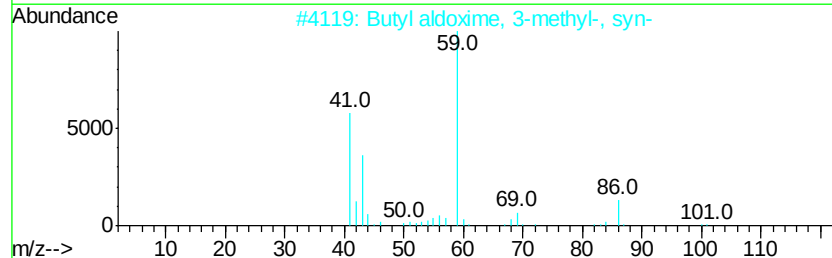
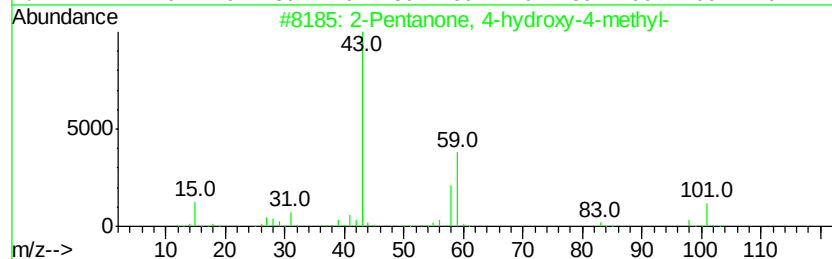
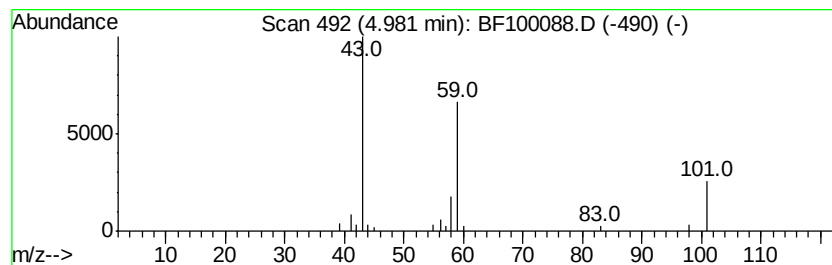
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown4.98 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	3.06 ng	73404	1,4-Dichlorobenzene-d4	6.77

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	42
2		Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	25
3		3-Octanol	130	C8H18O	000589-98-0	23
4		Guanidine	59	CH5N3	000113-00-8	9
5		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9



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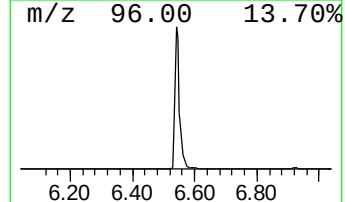
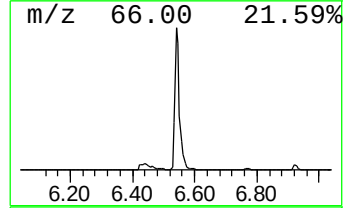
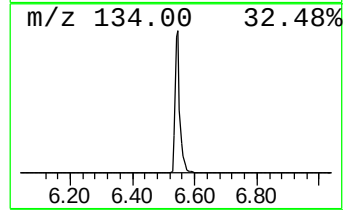
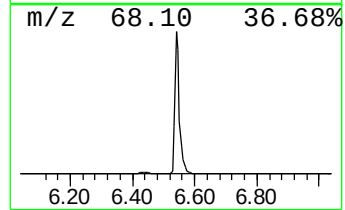
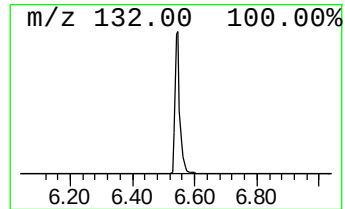
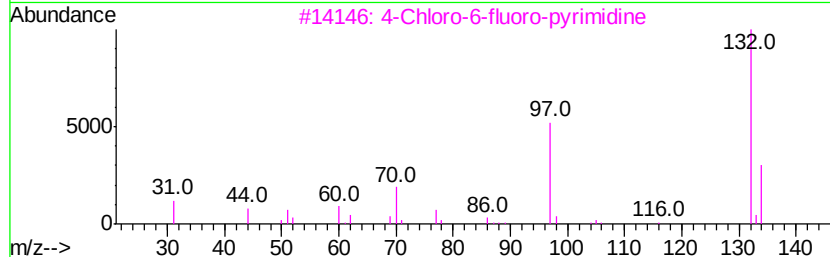
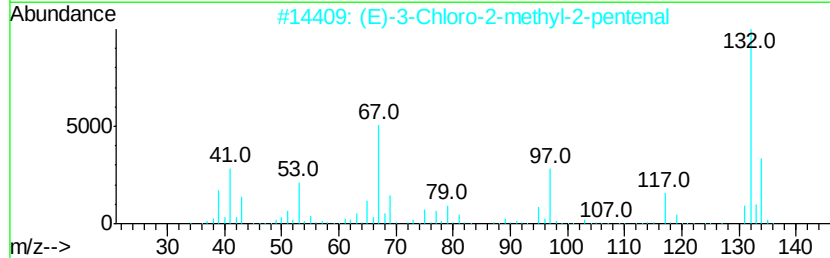
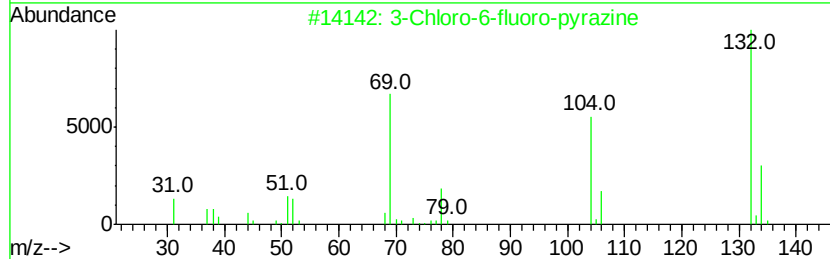
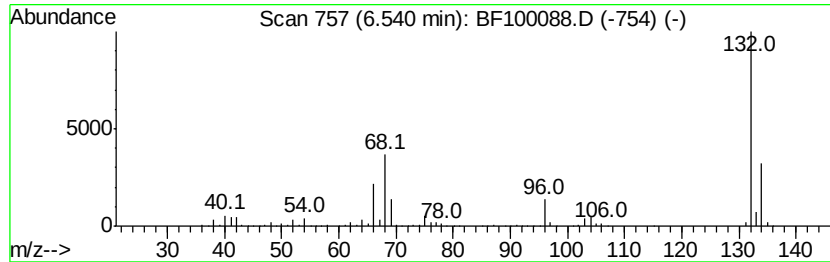
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Peak Number 2 unknown6.54 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.54	61.66 ng	1476950	1,4-Dichlorobenzene-d4	6.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
3		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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Peak Number 3 2-Propanol, 1-(2-methoxy-1-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	5.30 ng	126995	1,4-Dichlorobenzene-d4	6.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propanol, 1-(2-methoxy-1-methy...	148	C7H16O3	020324-32-7	72
2			2-Butanol, 3,3'-oxybis-	162	C8H18O3	054305-61-2	64
3			1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	42
4			2-Propanol, 1,1'-[(1-methyl-1,2-...	192	C9H20O4	001638-16-0	42
5			1-Propanol, 2-(2-methoxy-1-methy...	148	C7H16O3	055956-21-3	38

