

Data Path : Z:\HPCHEM1\BNA F\DATA\BF103017\
 Data File : BF099963.D
 Acq On : 30 Oct 2017 16:53
 Operator : SJ/JU
 Sample : I6171-01 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-102617-A

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	3.798	287	291	302	rBV	7169	13413	1.83%	0.219%
2	4.998	492	495	513	rBV	19299	35453	4.83%	0.578%
3	5.363	553	557	563	rBV2	10997	18000	2.45%	0.293%
4	5.416	563	566	582	rVV	127454	230520	31.39%	3.756%
5	5.622	598	601	605	rVB	8456	7885	1.07%	0.128%
6	6.434	737	739	758	rBV	179568	287464	39.15%	4.684%
7	6.563	758	761	767	rBV	303799	308648	42.03%	5.029%
8	6.786	796	799	814	rBB	497074	459508	62.58%	7.488%
9	6.939	822	825	834	rBV	252879	236356	32.19%	3.851%
10	7.357	894	896	911	rBV	131196	175664	23.92%	2.862%
11	8.069	1014	1017	1032	rBV	769769	651546	88.73%	10.617%
12	9.139	1196	1199	1209	rBV	514085	420831	57.31%	6.857%
13	9.545	1264	1268	1276	rVB	48219	45610	6.21%	0.743%
14	9.780	1303	1308	1312	rBV3	6890	11508	1.57%	0.188%
15	9.822	1312	1315	1328	rVB	813052	734295	100.00%	11.965%
16	10.539	1435	1437	1442	rBV	22279	20482	2.79%	0.334%
17	10.621	1448	1451	1464	rBV	190297	199295	27.14%	3.248%
18	10.721	1464	1468	1472	rVB2	5818	8348	1.14%	0.136%
19	11.321	1566	1570	1578	rBV	716641	723919	98.59%	11.796%
20	11.833	1654	1657	1660	rBV4	9609	8877	1.21%	0.145%
21	12.386	1747	1751	1757	rBV4	14587	26413	3.60%	0.430%
22	12.545	1775	1778	1784	rBV	12620	17999	2.45%	0.293%
23	12.627	1789	1792	1795	rBV4	7154	9573	1.30%	0.156%
24	12.710	1803	1806	1812	rVB7	16158	16856	2.30%	0.275%
25	12.762	1812	1815	1816	rBV3	20690	18994	2.59%	0.310%
26	12.874	1831	1834	1835	rBV3	9040	8556	1.17%	0.139%
27	12.904	1835	1839	1842	rVV	416329	398248	54.24%	6.489%
28	13.115	1873	1875	1878	rVB3	12403	9172	1.25%	0.149%
29	13.462	1931	1934	1936	rBV3	21101	21466	2.92%	0.350%
30	13.786	1987	1989	1992	rBV2	24991	23743	3.23%	0.387%
31	13.974	2018	2021	2029	rBV	507612	534601	72.80%	8.711%
32	14.421	2095	2097	2100	rBV3	23594	26399	3.60%	0.430%
33	15.468	2271	2275	2282	rVB	327472	427225	58.18%	6.962%

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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_F\METHODS\8270-BF102317.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

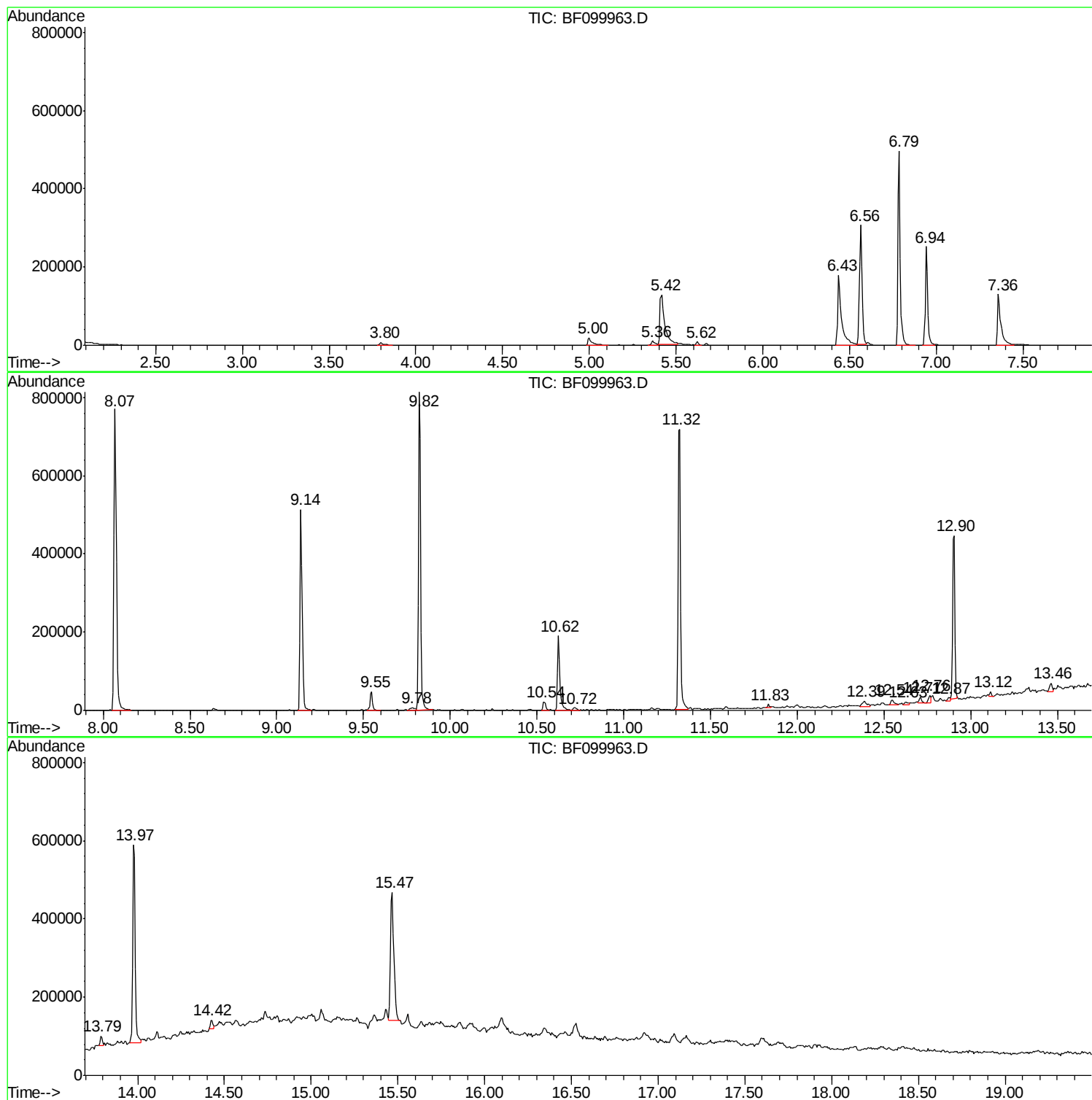
Sum of corrected areas: 6136867

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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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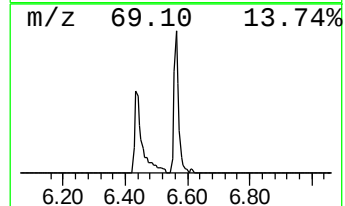
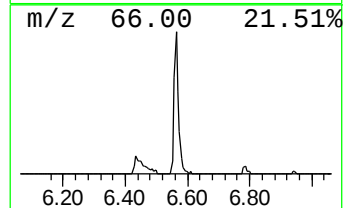
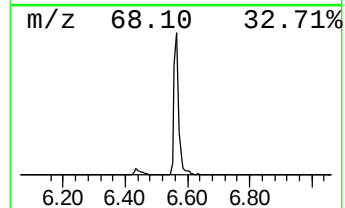
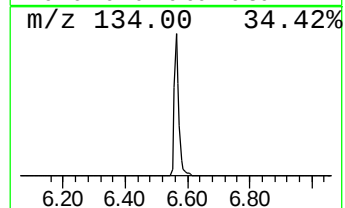
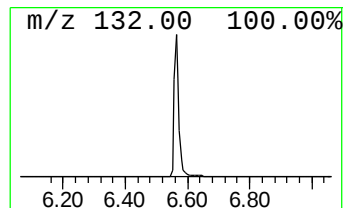
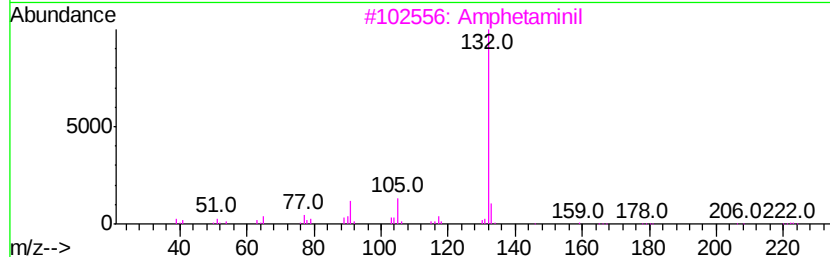
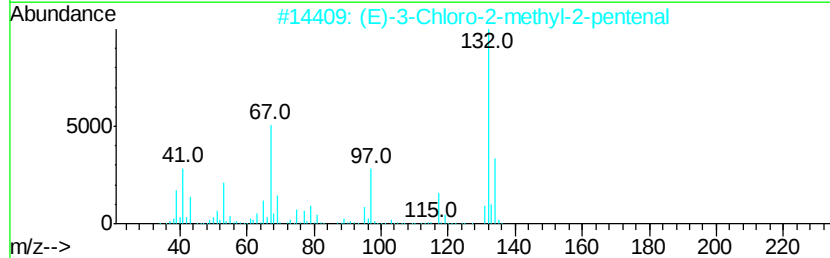
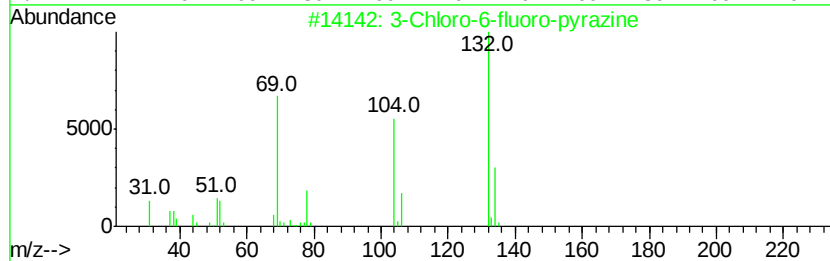
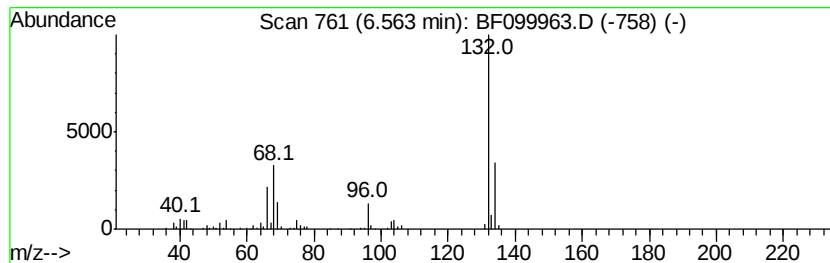
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.56	13.43 ng	308648	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Amphetaminil	250	C17H18N2	017590-01-1	12
4		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	12
5		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.56	6.56	13.4	ng	308648	1	6.79	459508	20.0