

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
 Data File : BF100154.D
 Acq On : 4 Nov 2017 00:08
 Operator : SJ/JU
 Sample : I6079-07 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q1-COMP

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	489	492	498	rBV	21648	33863	4.62%	0.515%
2	5.404	561	564	588	rBV	132432	302401	41.26%	4.602%
3	6.422	735	737	755	rBV	170941	347769	47.45%	5.293%
4	6.545	755	758	774	rVB	346470	416455	56.82%	6.338%
5	6.769	792	796	813	rVB	511905	517224	70.56%	7.872%
6	6.922	819	822	838	rBV	301311	313494	42.77%	4.771%
7	7.345	891	894	915	rBV	140316	215991	29.47%	3.287%
8	8.051	1011	1014	1032	rBV	752559	707112	96.47%	10.762%
9	9.122	1193	1196	1206	rBV	601750	471352	64.31%	7.174%
10	9.522	1261	1264	1274	rVB	88704	85580	11.68%	1.303%
11	9.804	1308	1312	1322	rVB	836282	732979	100.00%	11.156%
12	10.522	1432	1434	1439	rBB	16123	13270	1.81%	0.202%
13	10.604	1445	1448	1460	rBV	167542	178792	24.39%	2.721%
14	10.698	1460	1464	1469	rVB3	6291	8542	1.17%	0.130%
15	11.298	1562	1566	1576	rBV	703140	670855	91.52%	10.210%
16	11.839	1655	1658	1663	rVB	8127	8752	1.19%	0.133%
17	11.916	1669	1671	1674	rBV2	7929	8204	1.12%	0.125%
18	12.522	1771	1774	1779	rBV	51780	45008	6.14%	0.685%
19	12.733	1807	1810	1811	rBV2	9402	8838	1.21%	0.135%
20	12.751	1811	1813	1819	rVB	40835	39474	5.39%	0.601%
21	12.874	1831	1834	1838	rBV	357632	284663	38.84%	4.333%
22	13.086	1867	1870	1873	rVB	10405	10248	1.40%	0.156%
23	13.722	1972	1978	1981	rBV5	6029	9517	1.30%	0.145%
24	13.757	1981	1984	1991	rVV4	12702	17676	2.41%	0.269%
25	13.951	2012	2017	2021	rBV	430426	448502	61.19%	6.826%
26	14.669	2136	2139	2145	rBV2	33261	37180	5.07%	0.566%
27	14.998	2191	2195	2198	rBV2	31943	36565	4.99%	0.557%
28	15.298	2243	2246	2249	rBV	13287	13780	1.88%	0.210%
29	15.357	2252	2256	2260	rBV	32596	39649	5.41%	0.603%
30	15.410	2260	2265	2273	rVB	339396	431834	58.91%	6.572%
31	15.768	2323	2326	2332	rVB5	14285	22768	3.11%	0.347%
32	16.245	2402	2407	2414	rBV7	14487	30064	4.10%	0.458%
33	16.815	2498	2504	2507	rBV5	8269	15673	2.14%	0.239%
34	16.921	2517	2522	2526	rVB5	5715	9840	1.34%	0.150%

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35	17.374	2593	2599	2606	rVV7	5164	12790	1.74%	0.195%
36	17.474	2610	2616	2622	rVB6	11258	23683	3.23%	0.360%

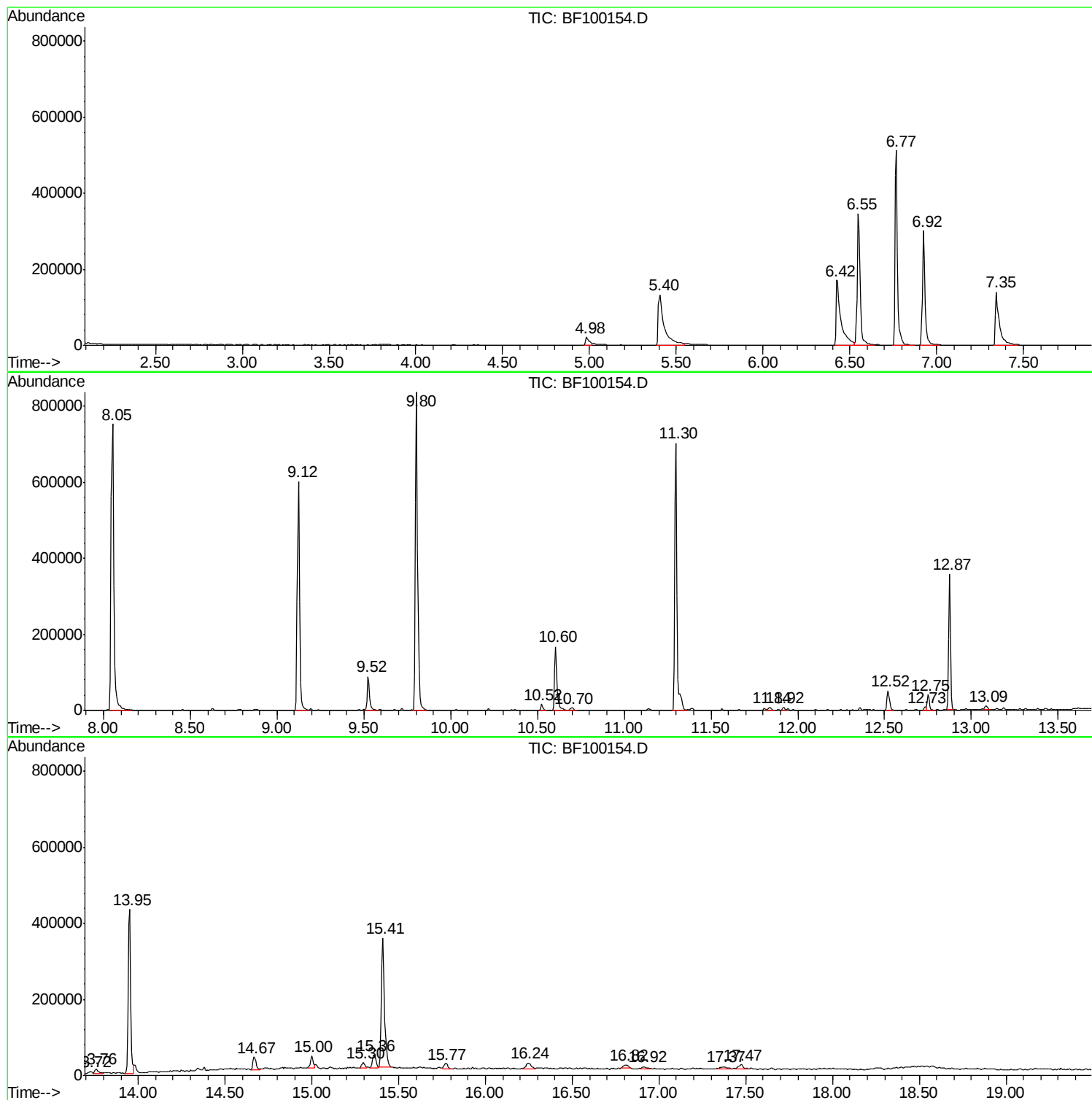
Sum of corrected areas: 6570387

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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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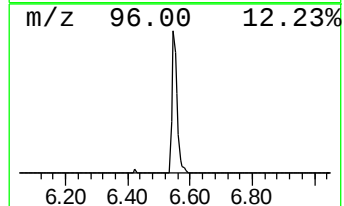
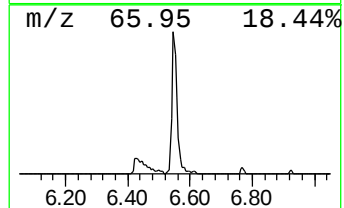
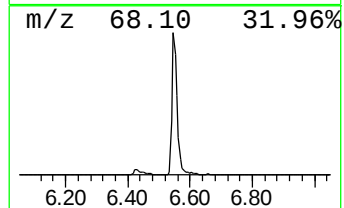
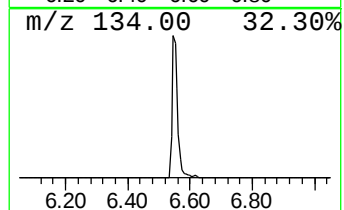
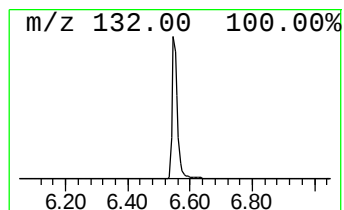
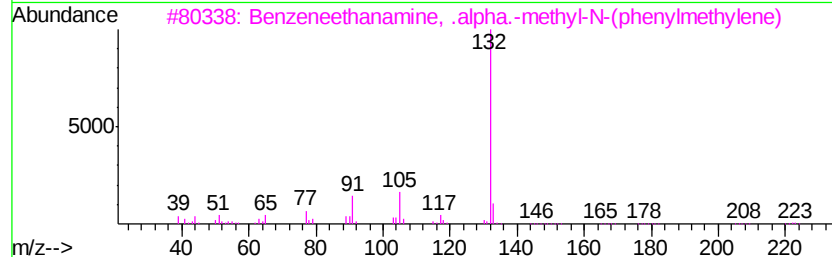
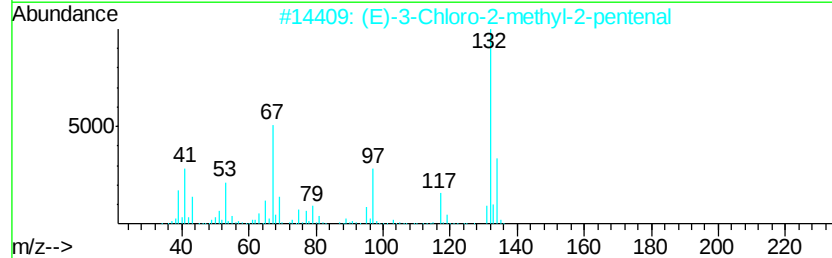
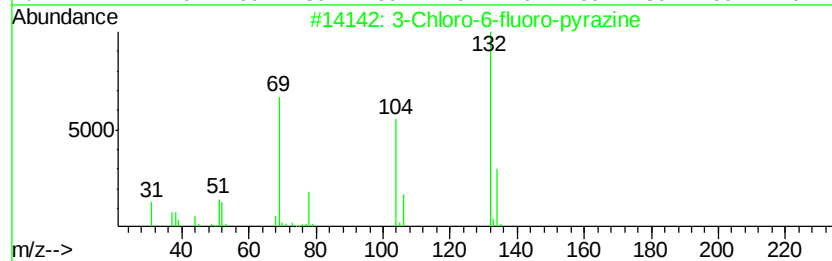
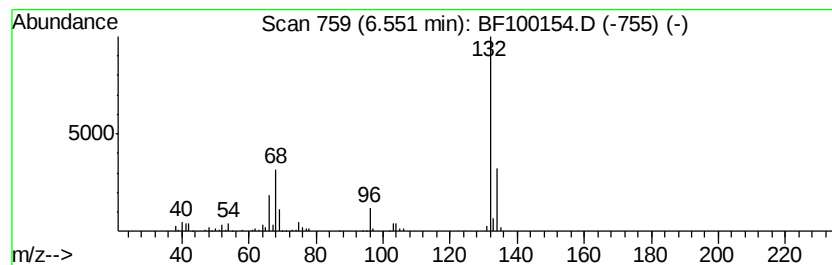
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Peak Number 1 unknown6.55 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.55	16.10 ng	416455	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	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Benzeneethanamine, .alpha.-methy...	223	C16H17N	002980-02-1	12
4		Amphetaminil	250	C17H18N2	017590-01-1	12
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.55	6.55	16.1	ng	416455	1	6.77	517224	20.0