

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF090717\
 Data File : BF098320.D
 Acq On : 7 Sep 2017 18:17
 Operator : SJ/JU
 Sample : I5052-01 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G26A-0-3-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-BF081517.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	4.851	465	470	474	rBV	17219	19510	1.96%	0.262%
2	5.281	540	543	560	rVB	226712	245061	24.68%	3.291%
3	6.322	717	720	733	rBV	272191	278802	28.08%	3.745%
4	6.451	739	742	746	rBV	312293	267467	26.93%	3.592%
5	6.686	778	782	789	rBV	945785	741806	74.70%	9.963%
6	6.839	805	808	812	rBV	243878	198919	20.03%	2.672%
7	7.251	874	878	882	rBV	169283	149263	15.03%	2.005%
8	7.969	996	1000	1004	rBV	1067415	978381	98.53%	13.141%
9	8.398	1069	1073	1077	rBV	12278	11112	1.12%	0.149%
10	8.533	1091	1096	1099	rBV	21722	19858	2.00%	0.267%
11	8.686	1119	1122	1125	rBV	21035	16003	1.61%	0.215%
12	8.780	1136	1138	1142	rBV	14595	11899	1.20%	0.160%
13	9.045	1179	1183	1187	rBV	379424	295374	29.74%	3.967%
14	9.133	1194	1198	1200	rBV	12213	13403	1.35%	0.180%
15	9.386	1238	1241	1243	rBV	11056	11000	1.11%	0.148%
16	9.445	1247	1251	1258	rVB2	39636	41770	4.21%	0.561%
17	9.586	1272	1275	1278	rBV	14102	12636	1.27%	0.170%
18	9.663	1285	1288	1292	rVB	13408	12477	1.26%	0.168%
19	9.721	1292	1298	1302	rBV	1068336	993021	100.00%	13.337%
20	9.927	1329	1333	1337	rVB	12932	12262	1.23%	0.165%
21	10.263	1388	1390	1398	rVB3	9388	10612	1.07%	0.143%
22	10.380	1405	1410	1413	rBV2	9363	11298	1.14%	0.152%
23	10.433	1416	1419	1423	rBV	66057	55432	5.58%	0.745%
24	10.510	1428	1432	1438	rVV	165383	135630	13.66%	1.822%
25	10.639	1449	1454	1458	rBV	28144	39967	4.02%	0.537%
26	11.080	1527	1529	1531	rVV2	14979	12345	1.24%	0.166%
27	11.110	1531	1534	1538	rVB3	12512	13821	1.39%	0.186%
28	11.204	1547	1550	1553	rBV	1007511	894638	90.09%	12.016%
29	11.227	1553	1554	1557	rVV	45125	26332	2.65%	0.354%
30	11.280	1561	1563	1566	rVB	14097	11952	1.20%	0.161%
31	11.733	1638	1640	1644	rVB3	15390	14301	1.44%	0.192%
32	11.815	1651	1654	1656	rBV	10652	10808	1.09%	0.145%
33	11.910	1667	1670	1672	rBV3	11001	10829	1.09%	0.145%
34	12.257	1726	1729	1731	rBV	15354	17801	1.79%	0.239%

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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_F\METHODS\8270-BF081517.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.415	1752	1756	1759	rBV	62991	55917	5.63%	0.751%
36	12.639	1789	1794	1797	rBV2	52787	62973	6.34%	0.846%
37	12.674	1797	1800	1802	rVV3	21257	22702	2.29%	0.305%
38	12.698	1802	1804	1808	rVB4	11429	14343	1.44%	0.193%
39	12.786	1816	1819	1827	rVB	282957	246327	24.81%	3.308%
40	12.857	1829	1831	1833	rBV3	12926	14176	1.43%	0.190%
41	12.968	1847	1850	1855	rVB6	18345	19995	2.01%	0.269%
42	13.274	1899	1902	1906	rBV2	27793	31130	3.13%	0.418%
43	13.839	1994	1998	2001	rBV	807550	785809	79.13%	10.554%
44	15.250	2234	2238	2246	rVB2	470320	596215	60.04%	8.008%

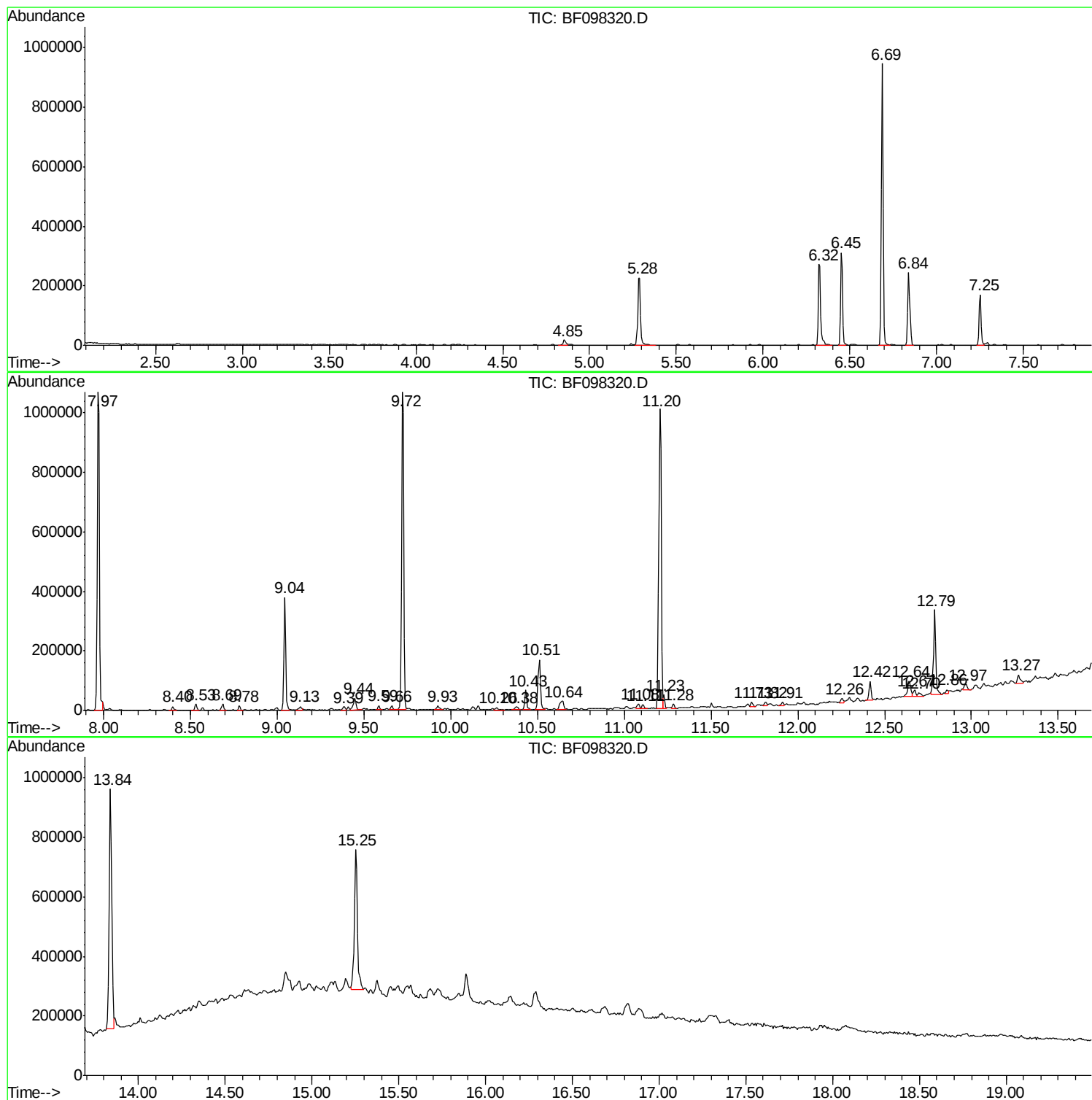
Sum of corrected areas: 7445377

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081517.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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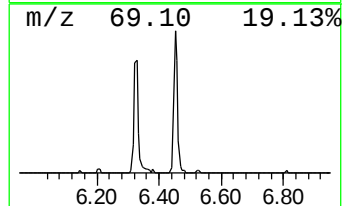
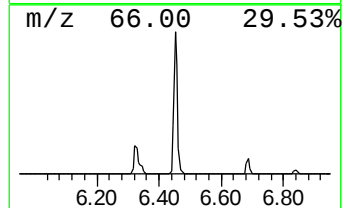
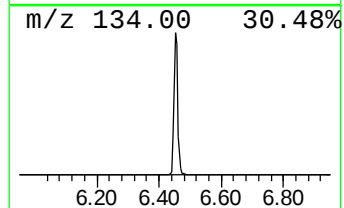
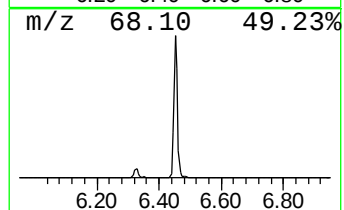
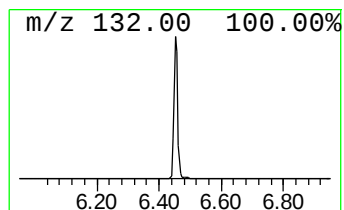
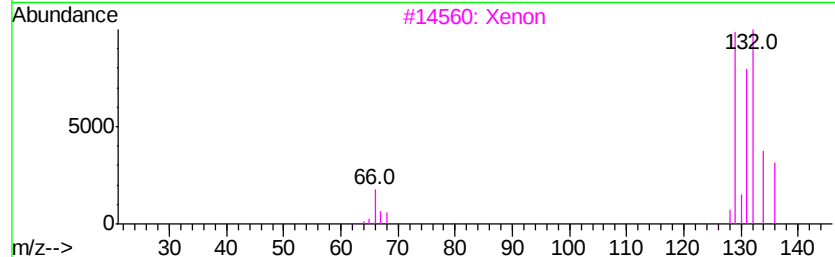
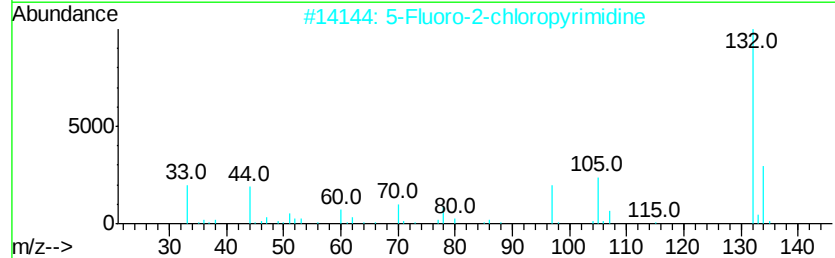
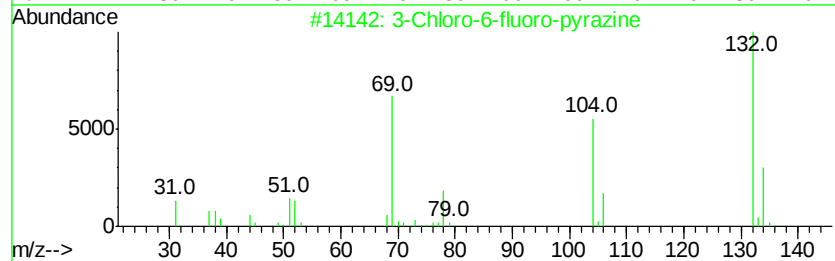
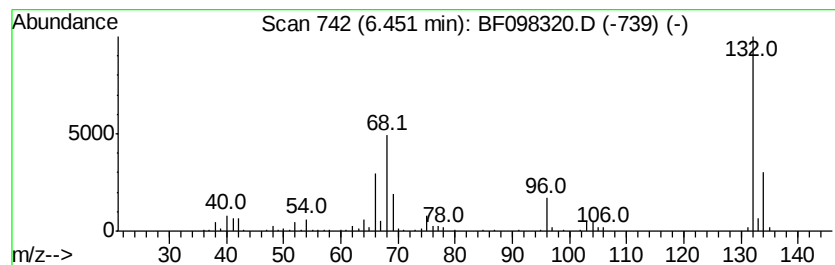
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.45	7.21 ng	267467	1,4-Dichlorobenzene-d4	6.69

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
2			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
3			Xenon	132	Xe	007440-63-3	9
4			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	9
5			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.45	6.45	7.2	ng	267467	1	6.69	741806	20.0