

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF090518\
 Data File : BF108789.D
 Acq On : 5 Sep 2018 22:07
 Operator : JU/SJ
 Sample : J4741-19 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AOC3-01-C

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF090518.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	5.325	505	508	512	rVB	415703	368802	13.67%	2.207%
2	6.348	679	682	686	rBV	483438	389741	14.45%	2.332%
3	6.495	703	707	710	rBV	498433	404824	15.01%	2.423%
4	6.731	744	747	751	rVB	2083848	1754869	65.06%	10.502%
5	6.889	770	774	777	rVB	395544	326565	12.11%	1.954%
6	7.284	837	841	846	rBV	301678	225719	8.37%	1.351%
7	8.013	961	965	969	rBV	2740433	2221442	82.35%	13.294%
8	9.089	1144	1148	1151	rBV	566831	446161	16.54%	2.670%
9	9.478	1211	1214	1218	rBV	116173	94225	3.49%	0.564%
10	9.766	1259	1263	1266	rBV	2311717	2131128	79.00%	12.754%
11	10.542	1391	1395	1399	rBV	173124	158894	5.89%	0.951%
12	11.236	1509	1513	1516	rBV	2131898	1879099	69.66%	11.246%
13	11.260	1516	1517	1520	rVB	174642	103033	3.82%	0.617%
14	11.313	1520	1526	1528	rBV	47692	42422	1.57%	0.254%
15	11.766	1600	1603	1608	rBV2	40351	61504	2.28%	0.368%
16	11.854	1614	1618	1623	rBV3	85953	125215	4.64%	0.749%
17	12.030	1645	1648	1650	rBV3	43850	47013	1.74%	0.281%
18	12.354	1700	1703	1704	rBV3	52347	57527	2.13%	0.344%
19	12.371	1704	1706	1710	rVB3	95372	86923	3.22%	0.520%
20	12.442	1715	1718	1722	rVB	463997	379799	14.08%	2.273%
21	12.548	1733	1736	1741	rBV3	94357	116752	4.33%	0.699%
22	12.589	1741	1743	1746	rVB3	39436	37978	1.41%	0.227%
23	12.666	1751	1756	1759	rBV	395514	448019	16.61%	2.681%
24	12.766	1771	1773	1775	rVB	46138	29356	1.09%	0.176%
25	12.818	1779	1782	1788	rVB	440460	454265	16.84%	2.719%
26	12.995	1810	1812	1817	rVB2	72100	101921	3.78%	0.610%
27	13.042	1818	1820	1822	rVV	75923	58315	2.16%	0.349%
28	13.066	1822	1824	1827	rVB	64734	54062	2.00%	0.324%
29	13.195	1844	1846	1849	rVB2	53261	45736	1.70%	0.274%
30	13.865	1955	1960	1963	rBV2	2505373	2697466	100.00%	16.143%
31	14.865	2128	2130	2133	rBV	169901	209409	7.76%	1.253%
32	15.271	2195	2199	2202	rVB	989830	1151376	42.68%	6.891%

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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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF090518.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

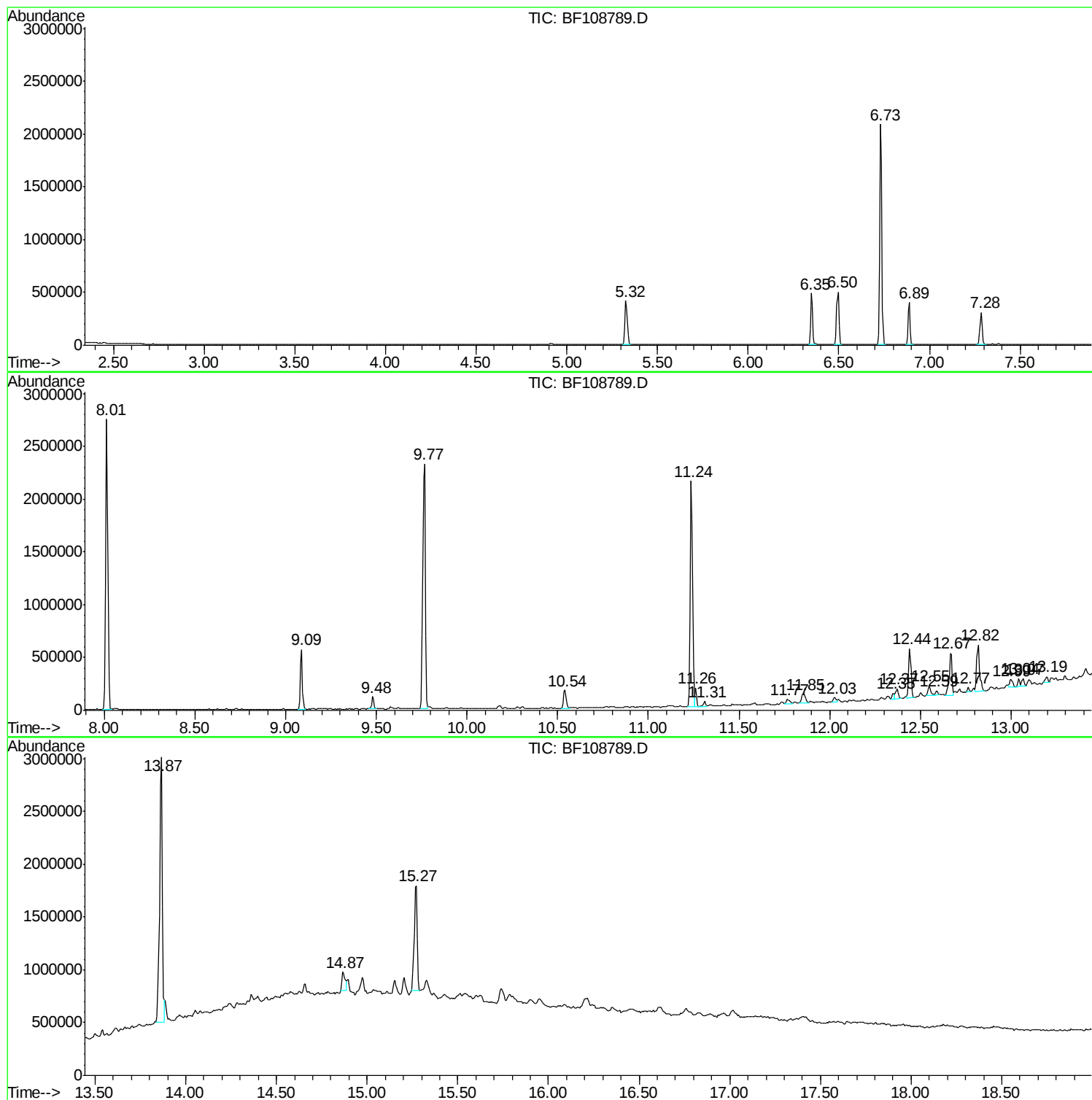
Sum of corrected areas: 16709560

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.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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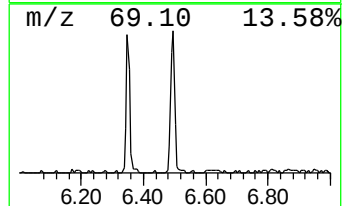
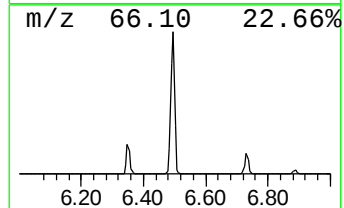
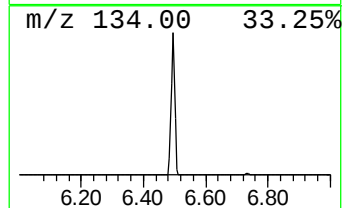
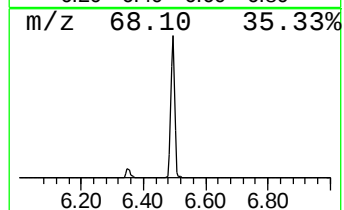
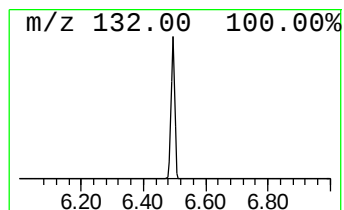
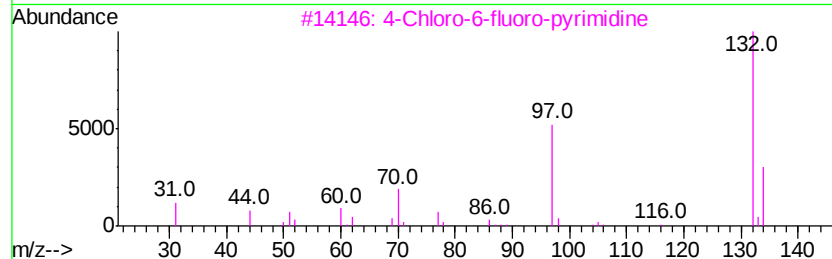
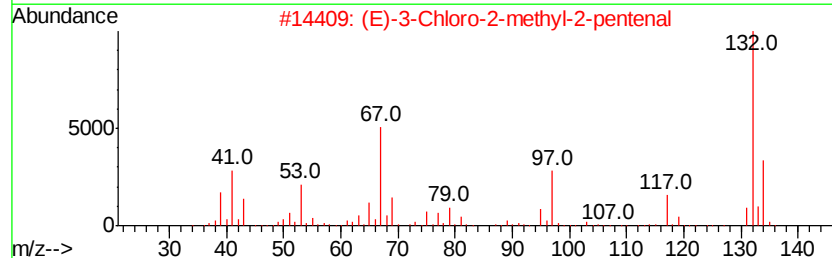
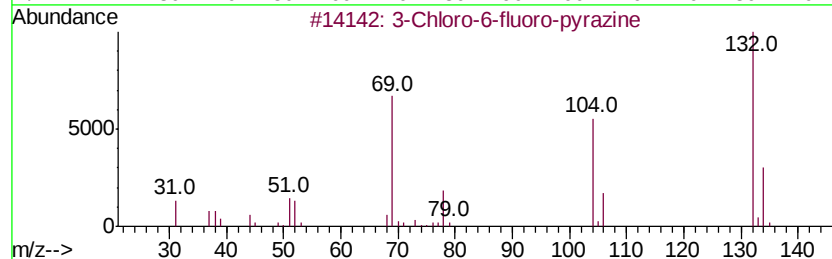
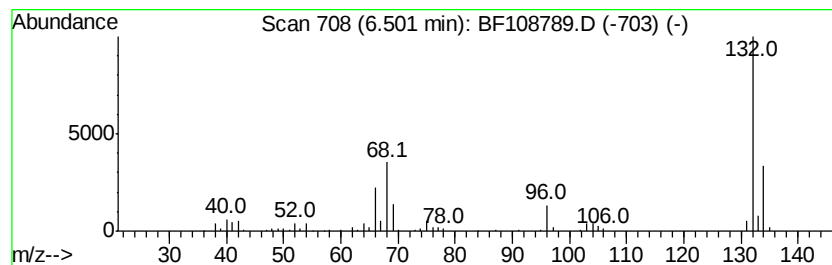
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Peak Number 1 unknown6.50 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.50	4.61 ng	404824	1,4-Dichlorobenzene-d4	6.73

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9
4			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5			4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.50	6.50	4.6	ng	404824	1	6.73	1754870	20.0