

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082218\
 Data File : BF108410.D
 Acq On : 23 Aug 2018 5:20
 Operator : JU/SJ
 Sample : J4525-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 081518-SIDI-F-D-S

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-BF080818.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	2.366	3	5	15	rVB	1344051	1578580	29.75%	5.417%
2	5.237	490	493	502	rBV	68863	75393	1.42%	0.259%
3	5.625	555	559	562	rBV	1555327	1355059	25.54%	4.650%
4	6.613	723	727	730	rBV	1088270	918294	17.31%	3.151%
5	6.643	730	732	740	rVB	113948	100452	1.89%	0.345%
6	6.772	749	754	757	rBV	2609438	2556496	48.18%	8.772%
7	7.001	789	793	796	rBV	1073254	915728	17.26%	3.142%
8	7.160	815	820	823	rBV	4045150	3609112	68.02%	12.384%
9	7.566	883	889	892	rBV	2758035	2995191	56.45%	10.278%
10	8.284	1007	1011	1023	rBV	1324263	1114492	21.00%	3.824%
11	9.360	1188	1194	1197	rBV	5834021	5306047	100.00%	18.207%
12	9.748	1256	1260	1265	rBV	110327	101780	1.92%	0.349%
13	10.042	1307	1310	1318	rVB	1234474	1156730	21.80%	3.969%
14	10.701	1419	1422	1425	rBV	124808	91571	1.73%	0.314%
15	10.836	1441	1445	1458	rBV	1605570	1450738	27.34%	4.978%
16	11.542	1561	1565	1568	rBV	1084204	995573	18.76%	3.416%
17	13.130	1830	1835	1838	rBV	3430772	3414200	64.35%	11.715%
18	14.189	2012	2015	2023	rBV	771436	770716	14.53%	2.645%
19	15.748	2273	2280	2288	rBV	446446	636876	12.00%	2.185%

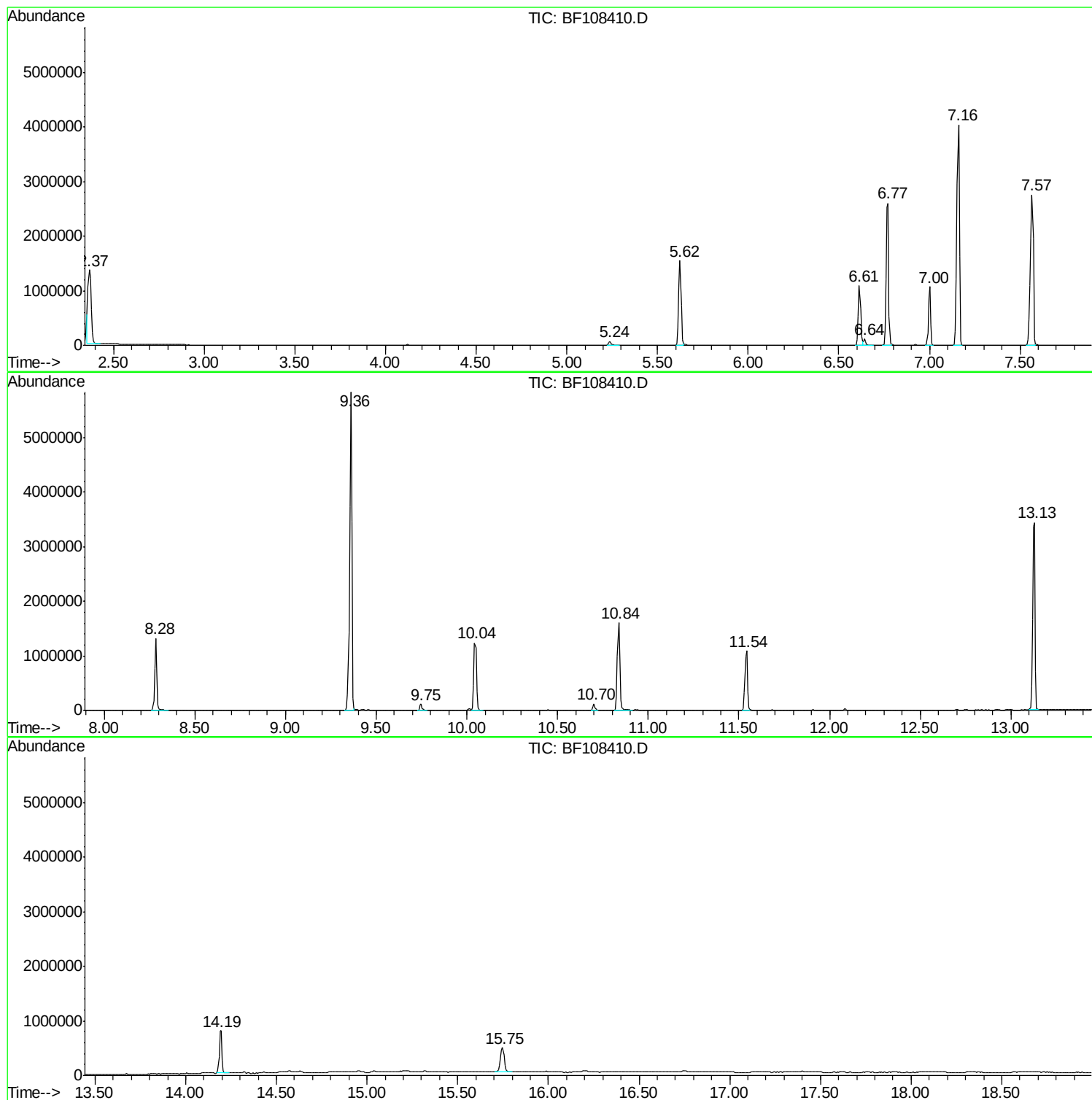
Sum of corrected areas: 29143028

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.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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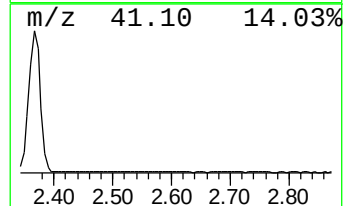
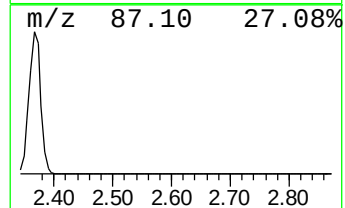
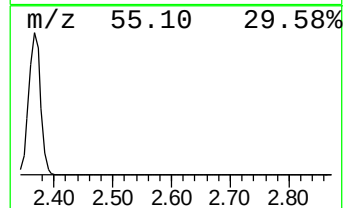
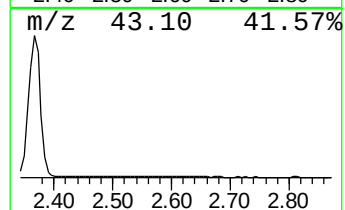
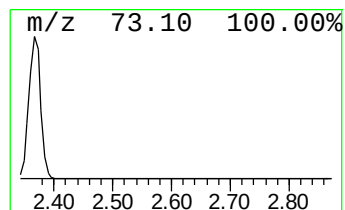
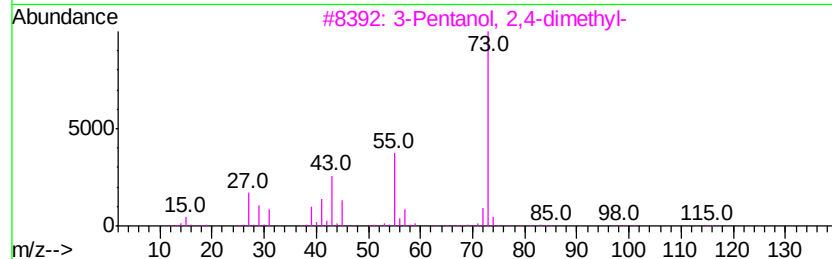
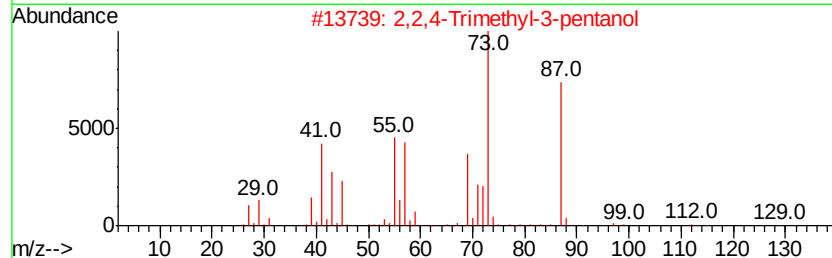
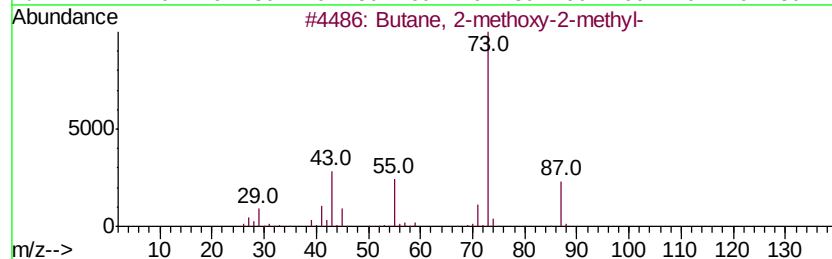
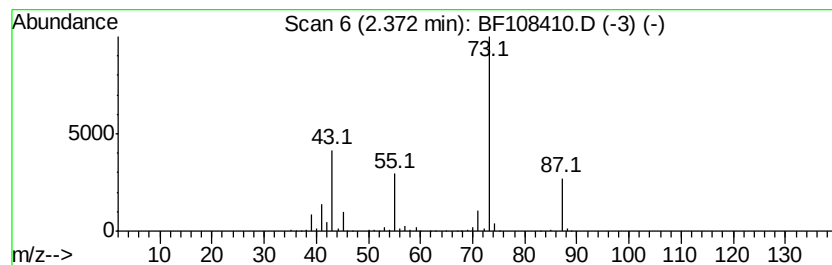
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.37	34.48 ng	1578580	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	38
3		3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	32
4		1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	28
5		Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17



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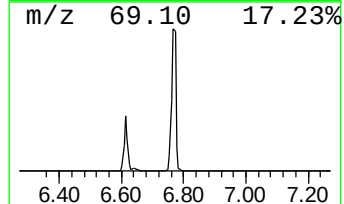
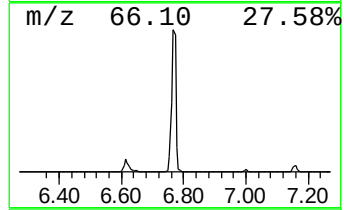
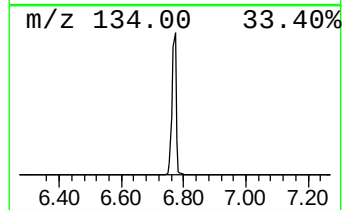
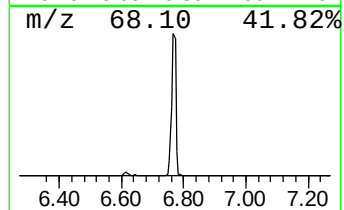
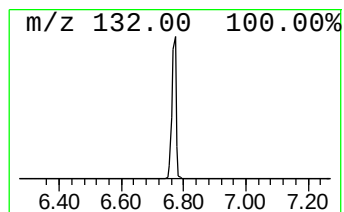
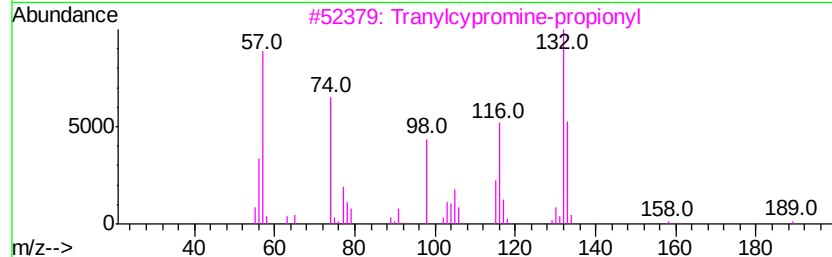
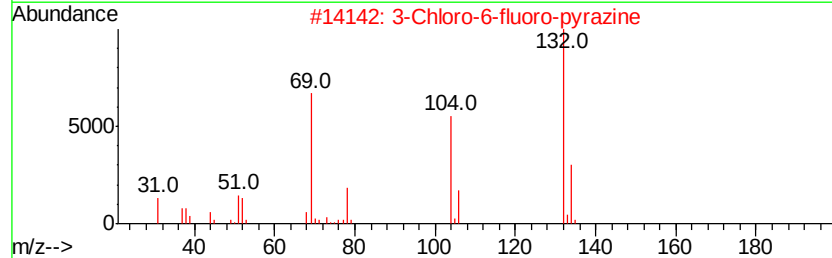
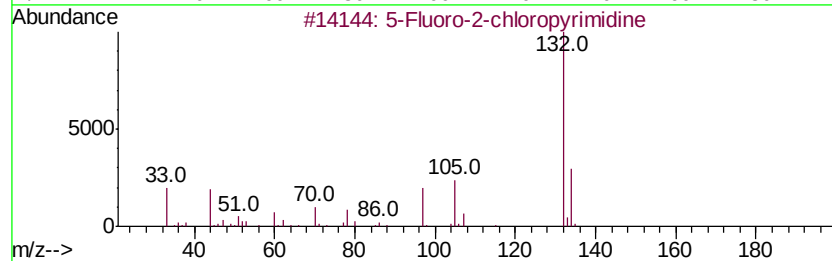
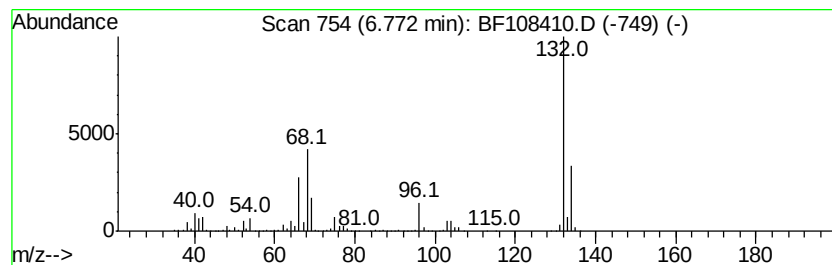
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Peak Number 2 unknown6.77 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.77	55.84 ng	2556500	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.37	34.5	ng	1578580	1	7.00	915728	20.0
unknown6.77	6.77	55.8	ng	2556500	1	7.00	915728	20.0