

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082718\
 Data File : BF108546.D
 Acq On : 28 Aug 2018 8:12
 Operator : JU/SJ
 Sample : J4660-02
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 082318-SIDI-F-D-M

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:\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	5.231	489	492	503	rVB	82342	91434	2.52%	0.455%
2	5.625	555	559	571	rBV	1150543	1049540	28.92%	5.220%
3	6.613	724	727	744	rBV	784524	699727	19.28%	3.480%
4	6.766	749	753	757	rBV	2118997	1858275	51.21%	9.242%
5	6.995	789	792	796	rBV	693491	633113	17.45%	3.149%
6	7.154	815	819	823	rBV	2829803	2616652	72.11%	13.014%
7	7.566	883	889	892	rBV	2037586	2104147	57.98%	10.465%
8	8.284	1007	1011	1026	rBV	778719	722064	19.90%	3.591%
9	9.354	1189	1193	1197	rBV	3828412	3628865	100.00%	18.049%
10	9.742	1256	1259	1265	rBV	58621	58703	1.62%	0.292%
11	10.042	1307	1310	1323	rVB	773945	676860	18.65%	3.366%
12	10.701	1419	1422	1425	rVB	51784	40162	1.11%	0.200%
13	10.836	1441	1445	1458	rBV	1211884	1144224	31.53%	5.691%
14	11.536	1561	1564	1573	rBV	663675	593149	16.35%	2.950%
15	13.124	1829	1834	1837	rBV	3284293	2937464	80.95%	14.610%
16	14.189	2011	2015	2026	rVB	769572	709822	19.56%	3.530%
17	15.036	2155	2159	2163	rVB	36749	40302	1.11%	0.200%
18	15.748	2272	2280	2288	rVB	335286	501309	13.81%	2.493%

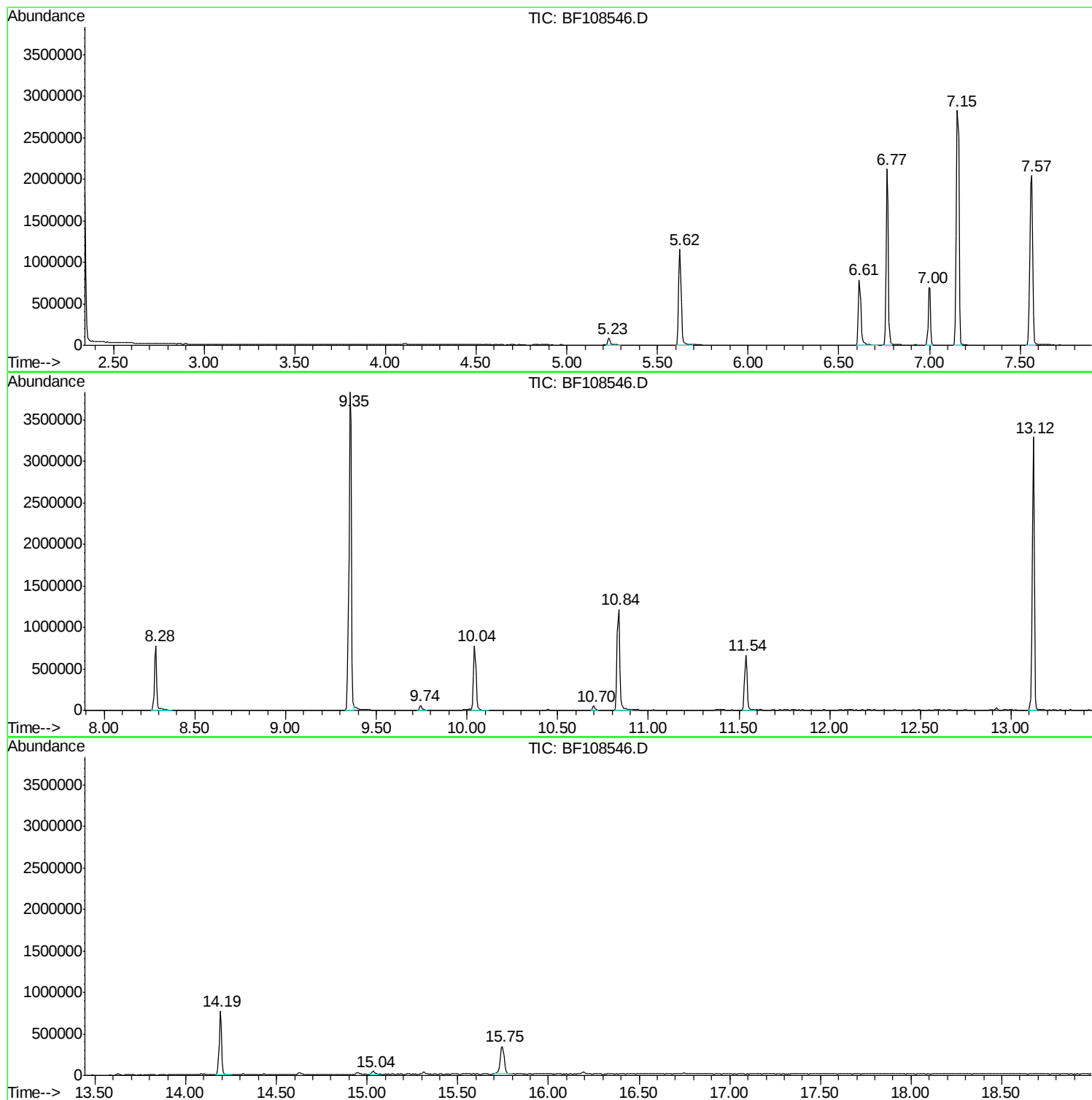
Sum of corrected areas: 20105812

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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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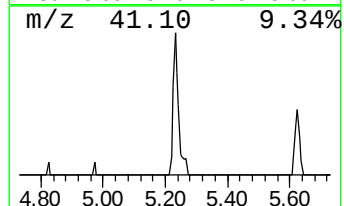
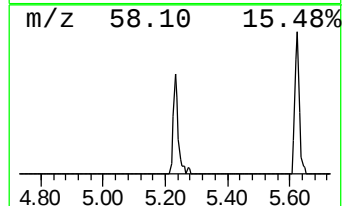
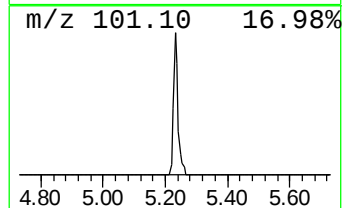
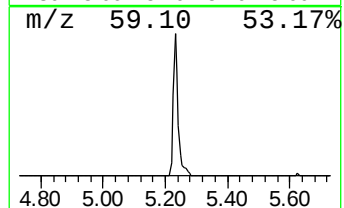
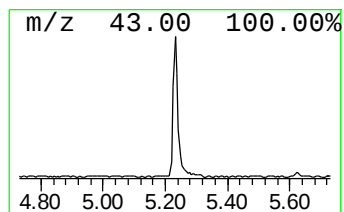
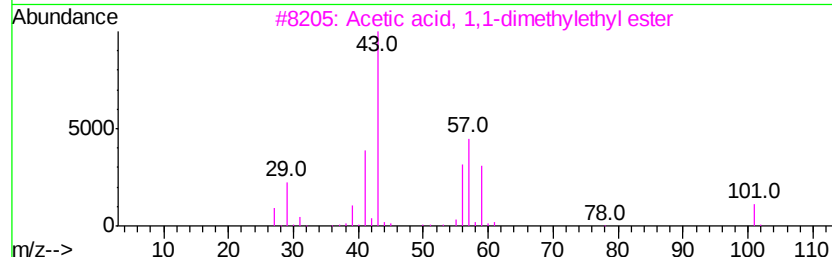
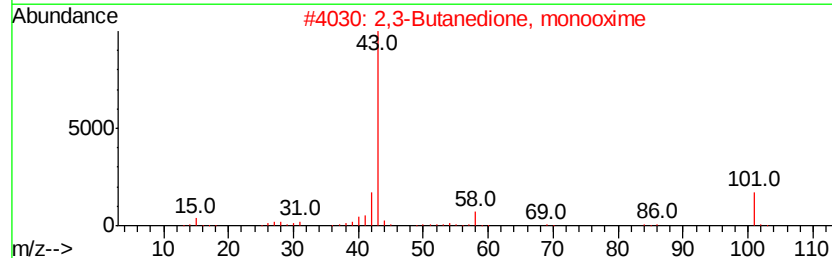
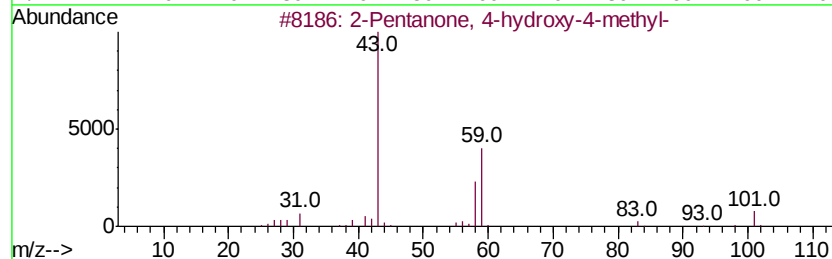
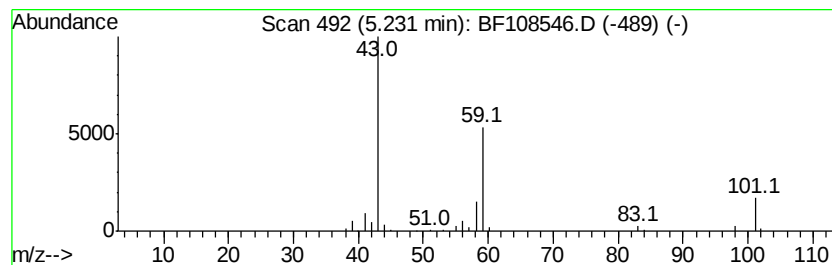
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TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.23	2.89 ng	91434	1,4-Dichlorobenzene-d4	7.00

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	25
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	23
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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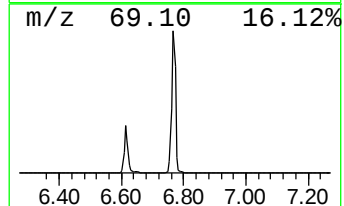
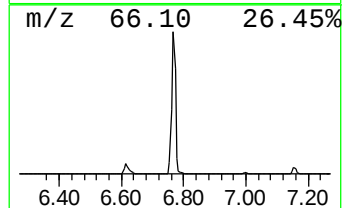
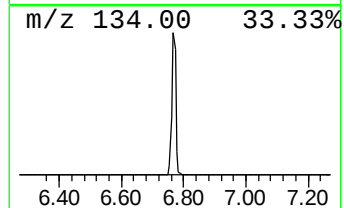
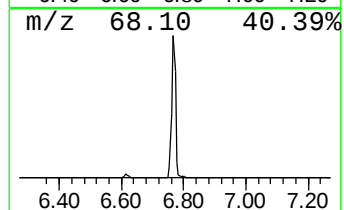
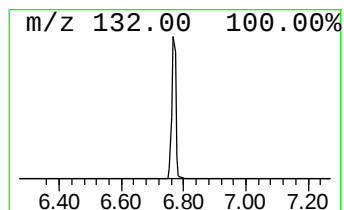
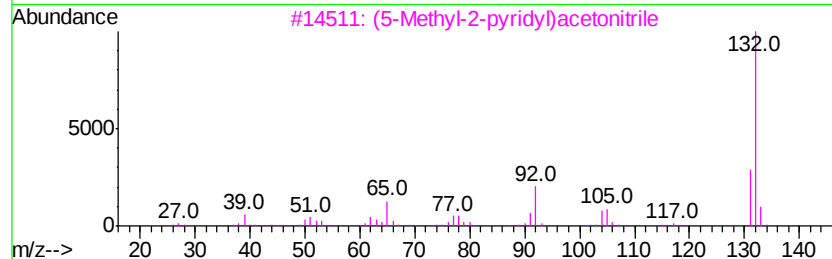
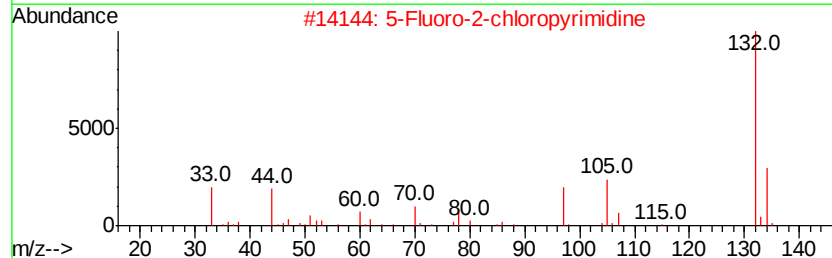
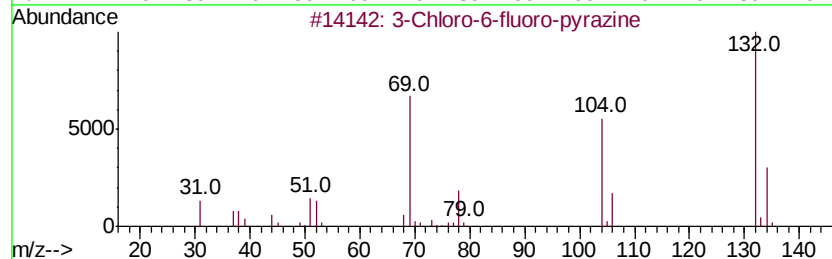
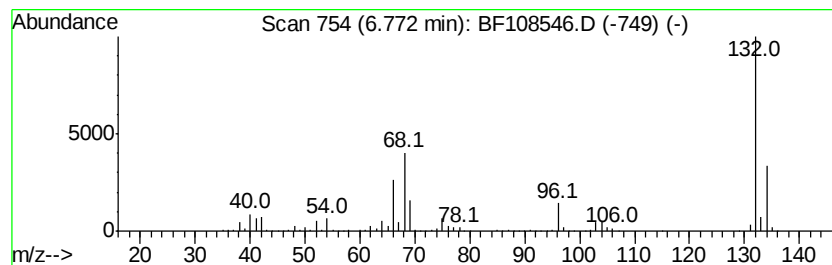
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Peak Number 2 unknown6.77 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.77	58.70 ng	1858280	1,4-Dichlorobenzene-d4	7.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.23	2.9	ng	91434	1	7.00	633113	20.0
unknown6.77	6.77	58.7	ng	1858280	1	7.00	633113	20.0