

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF082218\
 Data File : BF108411.D
 Acq On : 23 Aug 2018 5:47
 Operator : JU/SJ
 Sample : J4525-12
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 081518-SIDI-F-U-B

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	2.366	3	5	16	rVB	1323195	1365694	26.70%	4.855%
2	5.237	489	493	503	rVB	63185	73881	1.44%	0.263%
3	5.625	555	559	570	rBV	1429538	1251538	24.47%	4.449%
4	6.613	723	727	730	rBV	978114	815861	15.95%	2.901%
5	6.642	730	732	741	rVB	106124	96610	1.89%	0.343%
6	6.772	749	754	757	rBV	2527263	2429400	47.49%	8.637%
7	7.001	789	793	796	rBV	1095211	902365	17.64%	3.208%
8	7.160	815	820	823	rBV	3860222	3475732	67.94%	12.357%
9	7.566	883	889	892	rBV	2802868	2893793	56.57%	10.288%
10	8.283	1007	1011	1014	rBV	1344984	1079578	21.10%	3.838%
11	9.360	1188	1194	1197	rBV	5708029	5115591	100.00%	18.187%
12	9.748	1256	1260	1265	rBV	180587	153334	3.00%	0.545%
13	10.048	1307	1311	1314	rBV	1241159	1149109	22.46%	4.085%
14	10.701	1419	1422	1425	rBV	142449	103177	2.02%	0.367%
15	10.836	1441	1445	1454	rBV	1609160	1433623	28.02%	5.097%
16	11.542	1561	1565	1568	rBV	1162221	1012585	19.79%	3.600%
17	13.124	1830	1834	1838	rBV	3298427	3320481	64.91%	11.805%
18	14.189	2012	2015	2019	rBV	792135	768011	15.01%	2.730%
19	15.742	2273	2279	2288	rBV	465238	687378	13.44%	2.444%

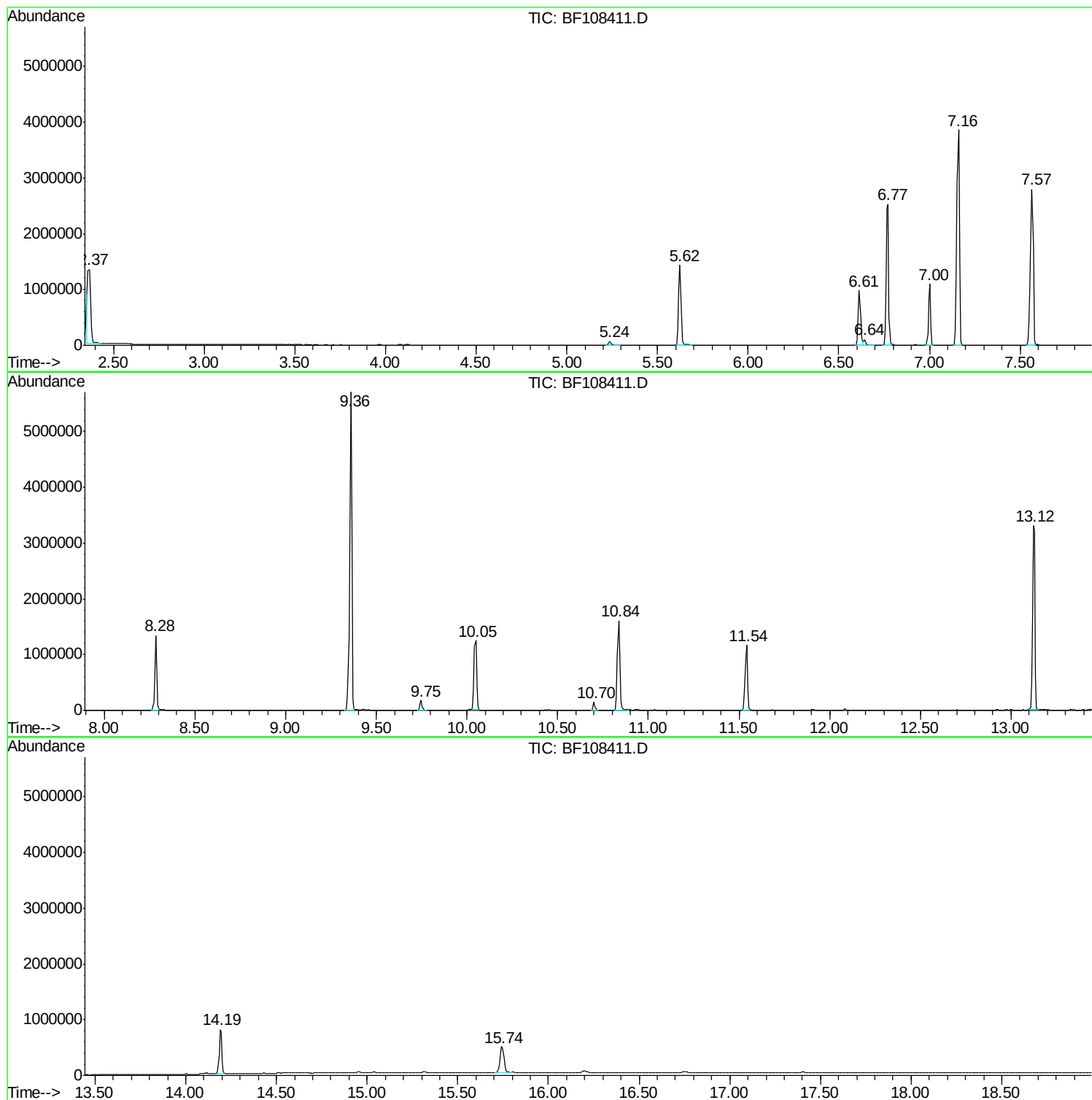
Sum of corrected areas: 28127741

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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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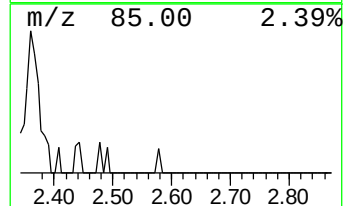
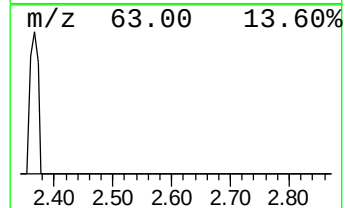
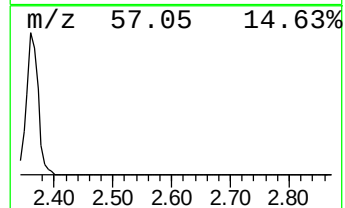
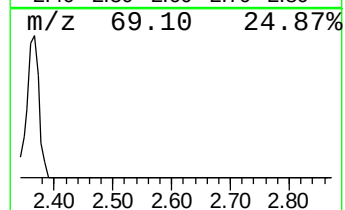
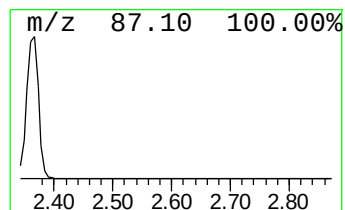
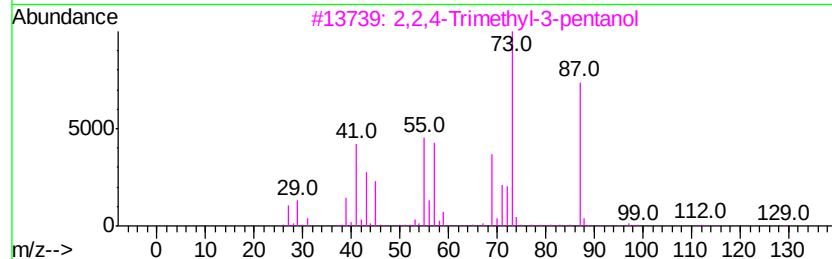
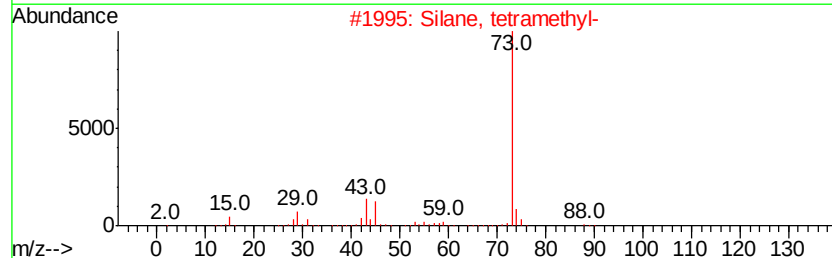
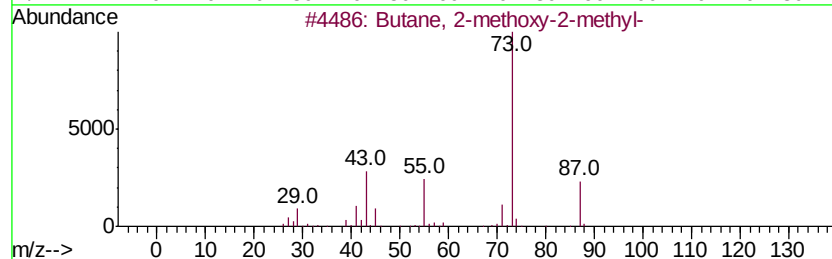
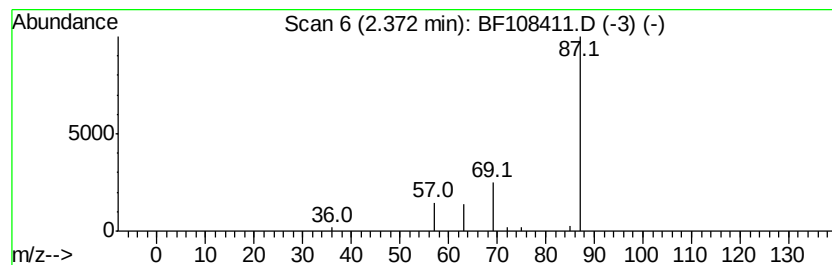
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.37	30.27 ng	1365690	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	78
2		Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3		2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
4		2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	25
5		Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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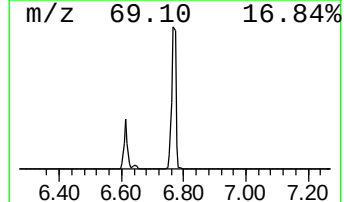
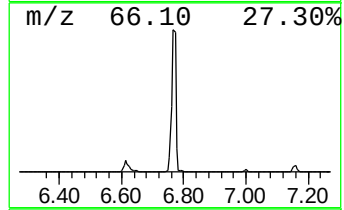
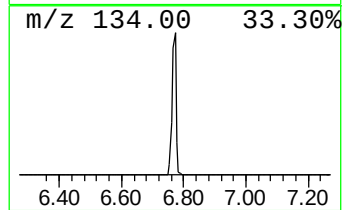
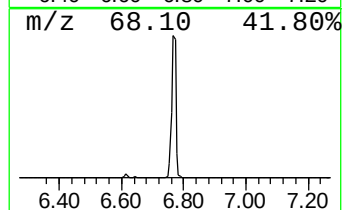
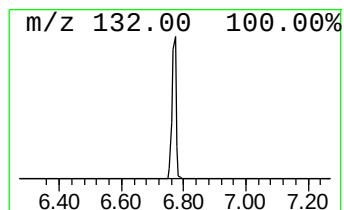
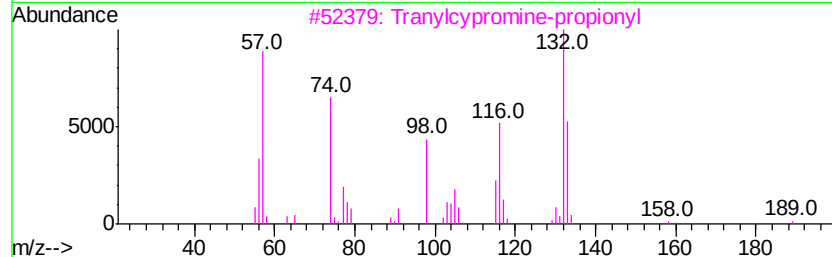
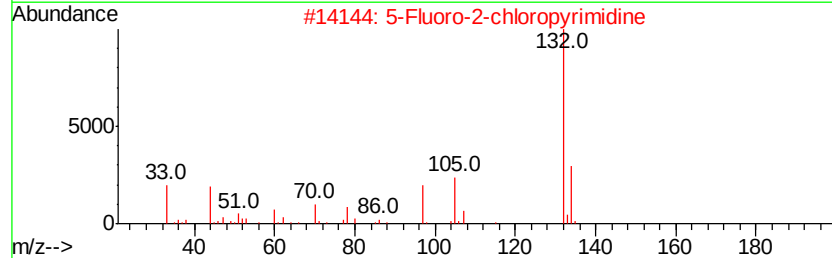
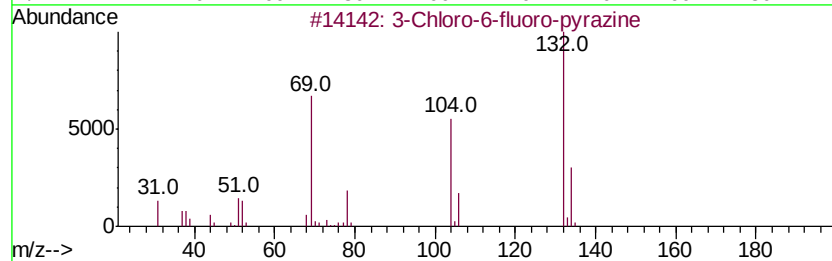
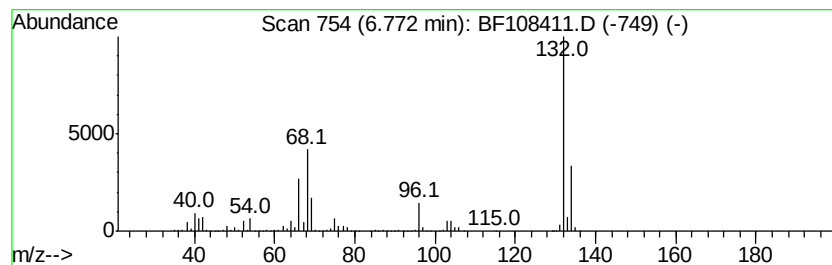
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Peak Number 2 unknown6.77 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.77	53.85 ng	2429400	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	35
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	16
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	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	30.3	ng	1365690	1	7.00	902365	20.0
unknown6.77	6.77	53.9	ng	2429400	1	7.00	902365	20.0