

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF110818\
 Data File : BF110613.D
 Acq On : 9 Nov 2018 4:59
 Operator : JU/SJ
 Sample : PB114627BL
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB114627BL

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-BF110818.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	5.187	523	527	537	rBV	70795	97170	3.72%	0.615%
2	5.569	588	592	612	rBV	1624068	1533330	58.65%	9.702%
3	6.569	757	762	782	rBV	1613216	1542850	59.02%	9.763%
4	6.710	782	786	790	rBV	1697525	1605116	61.40%	10.157%
5	6.945	822	826	836	rBB	482408	450329	17.23%	2.850%
6	7.098	848	852	856	rBV	1975830	1821662	69.68%	11.527%
7	7.510	917	922	925	rBV	1458271	1399475	53.53%	8.855%
8	8.222	1039	1043	1061	rBV	547815	597210	22.84%	3.779%
9	9.298	1221	1226	1229	rBV	2819573	2614248	100.00%	16.542%
10	9.986	1338	1343	1354	rBV	594697	588370	22.51%	3.723%
11	10.775	1473	1477	1490	rBV	753507	701091	26.82%	4.436%
12	11.475	1593	1596	1614	rBB	443039	474054	18.13%	3.000%
13	13.057	1861	1865	1869	rBV	1632307	1558146	59.60%	9.859%
14	14.127	2043	2047	2057	rBV	456814	431796	16.52%	2.732%
15	15.227	2227	2234	2242	rBV2	18758	29990	1.15%	0.190%
16	15.651	2302	2306	2314	rVB	248989	328693	12.57%	2.080%
17	16.074	2373	2378	2383	rVB2	19896	30114	1.15%	0.191%

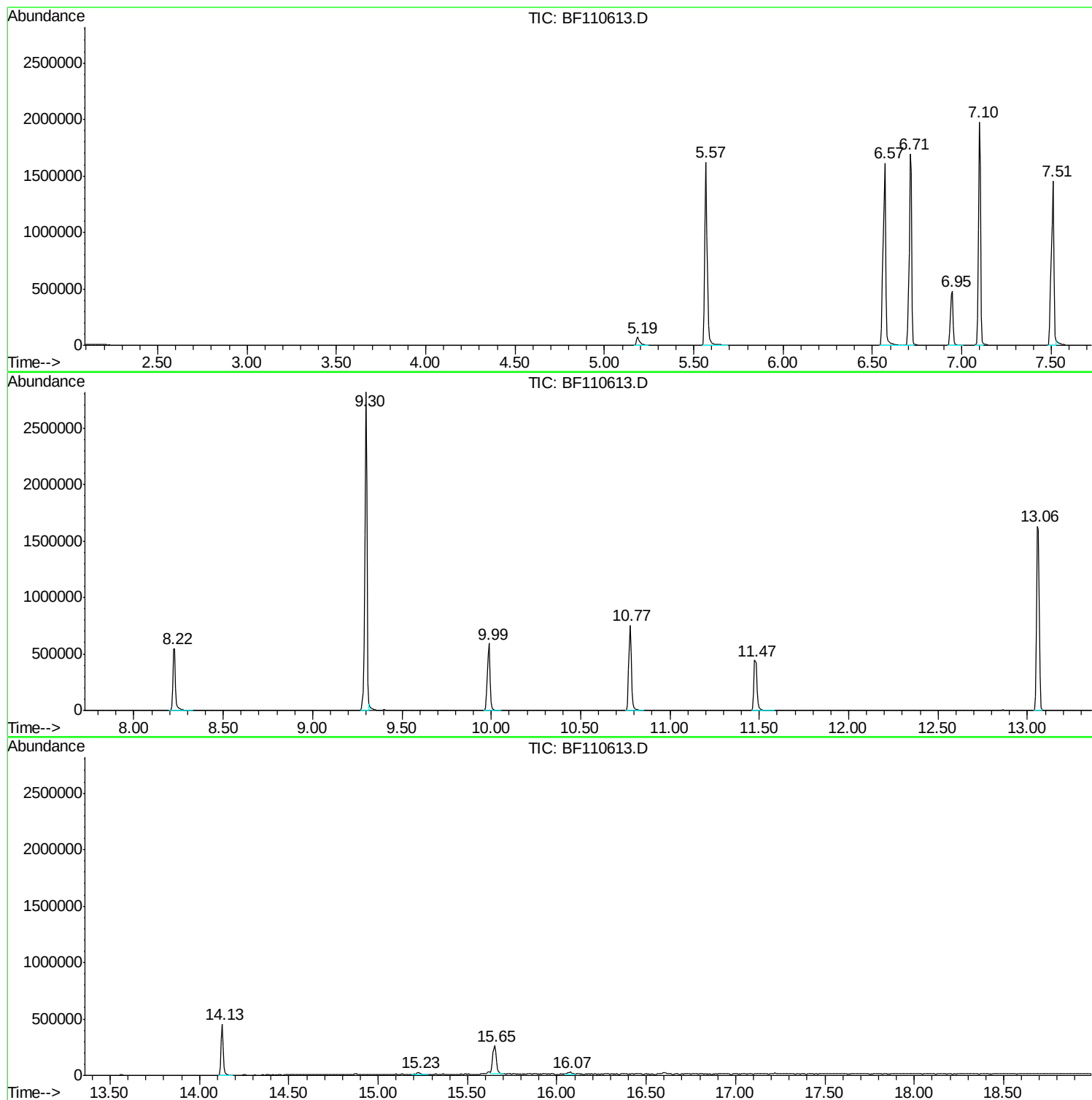
Sum of corrected areas: 15803644

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.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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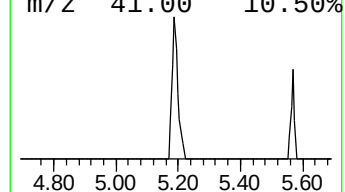
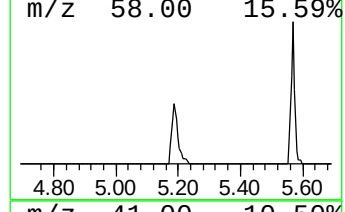
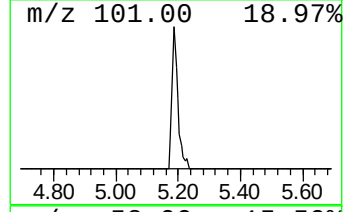
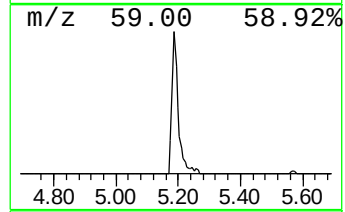
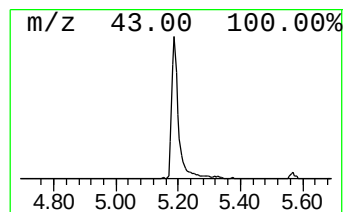
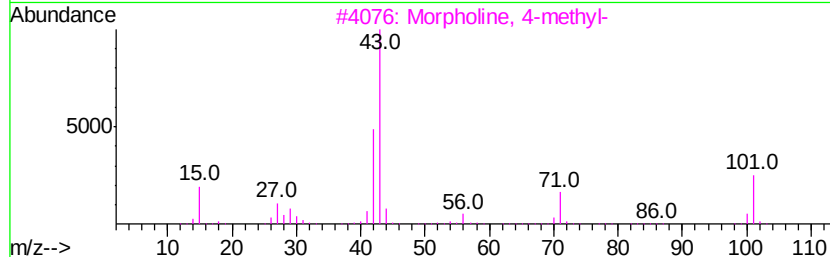
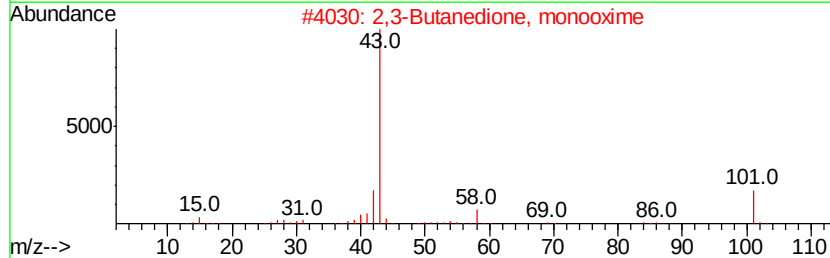
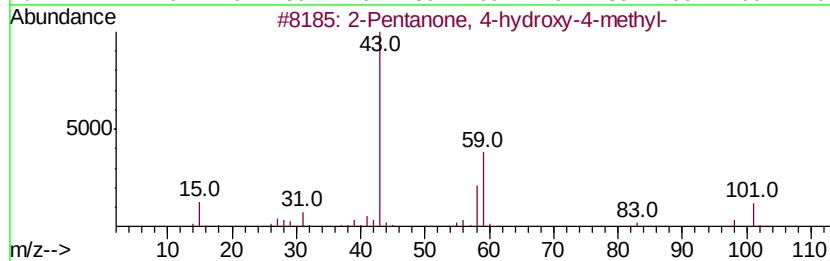
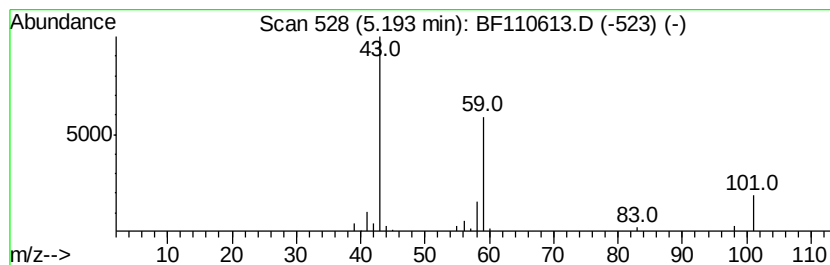
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	4.32 ng	97170	1,4-Dichlorobenzene-d4	6.95

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	35
3			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	9
5			Propanoic acid, 2-nitro-, methyl...	133	C4H7NO4	006118-50-9	9



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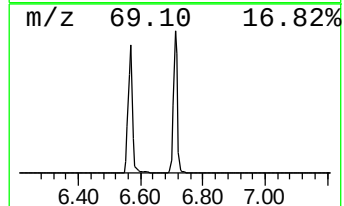
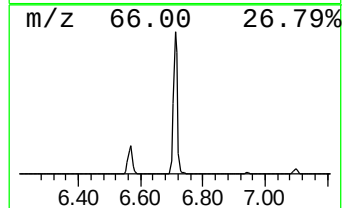
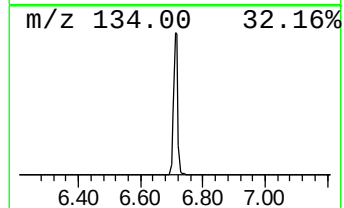
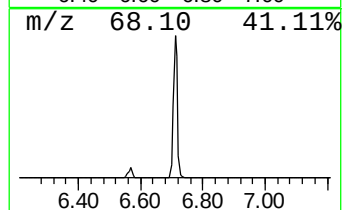
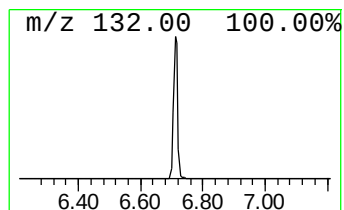
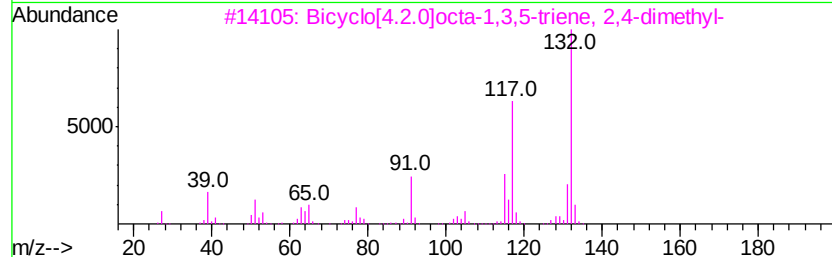
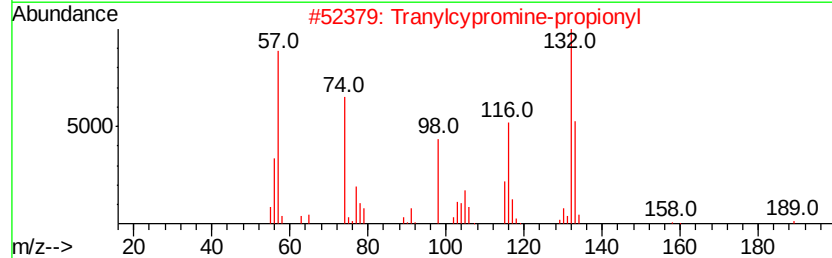
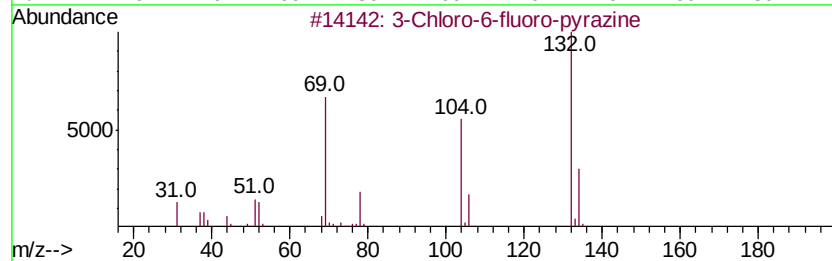
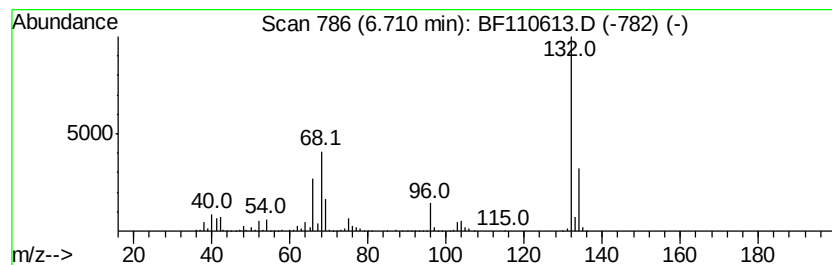
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Peak Number 2 unknown6.71 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.71	71.29 ng	1605120	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		5-Aminoindole	132	C8H8N2	005192-03-0	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	4.3	ng	97170	1	6.95	450329	20.0
unknown6.71	6.71	71.3	ng	1605120	1	6.95	450329	20.0