

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF080919\
 Data File : BF116106.D
 Acq On : 9 Aug 2019 11:57
 Operator : HP/JU
 Sample : PB122075BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122075BL

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-BF073019.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	2.193	14	18	29	rVB	164843	243492	5.65%	0.719%
2	5.475	572	576	600	rBV	3243357	3427892	79.49%	10.117%
3	6.481	742	747	765	rBV	3169135	3573521	82.86%	10.547%
4	6.616	765	770	773	rBV	3555322	3536471	82.00%	10.438%
5	6.839	804	808	820	rVB	934055	851768	19.75%	2.514%
6	6.992	830	834	849	rBV	2910671	2965175	68.76%	8.751%
7	7.404	899	904	928	rBV	2025487	2437699	56.53%	7.195%
8	8.116	1021	1025	1049	rBV	997004	1096975	25.44%	3.238%
9	9.192	1203	1208	1211	rBV	4296679	4189561	97.15%	12.365%
10	9.869	1319	1323	1344	rBV	1192679	1285011	29.80%	3.793%
11	10.663	1454	1458	1488	rBV	2444303	2671031	61.94%	7.883%
12	11.357	1572	1576	1601	rBV	1188595	1280774	29.70%	3.780%
13	12.945	1841	1846	1849	rBV	3927306	4312577	100.00%	12.728%
14	13.998	2021	2025	2043	rBV	813966	1007908	23.37%	2.975%
15	15.468	2270	2275	2296	rVB	530581	880951	20.43%	2.600%
16	15.892	2342	2347	2356	rVB	42258	68310	1.58%	0.202%
17	16.386	2426	2431	2437	rBV2	30281	53000	1.23%	0.156%

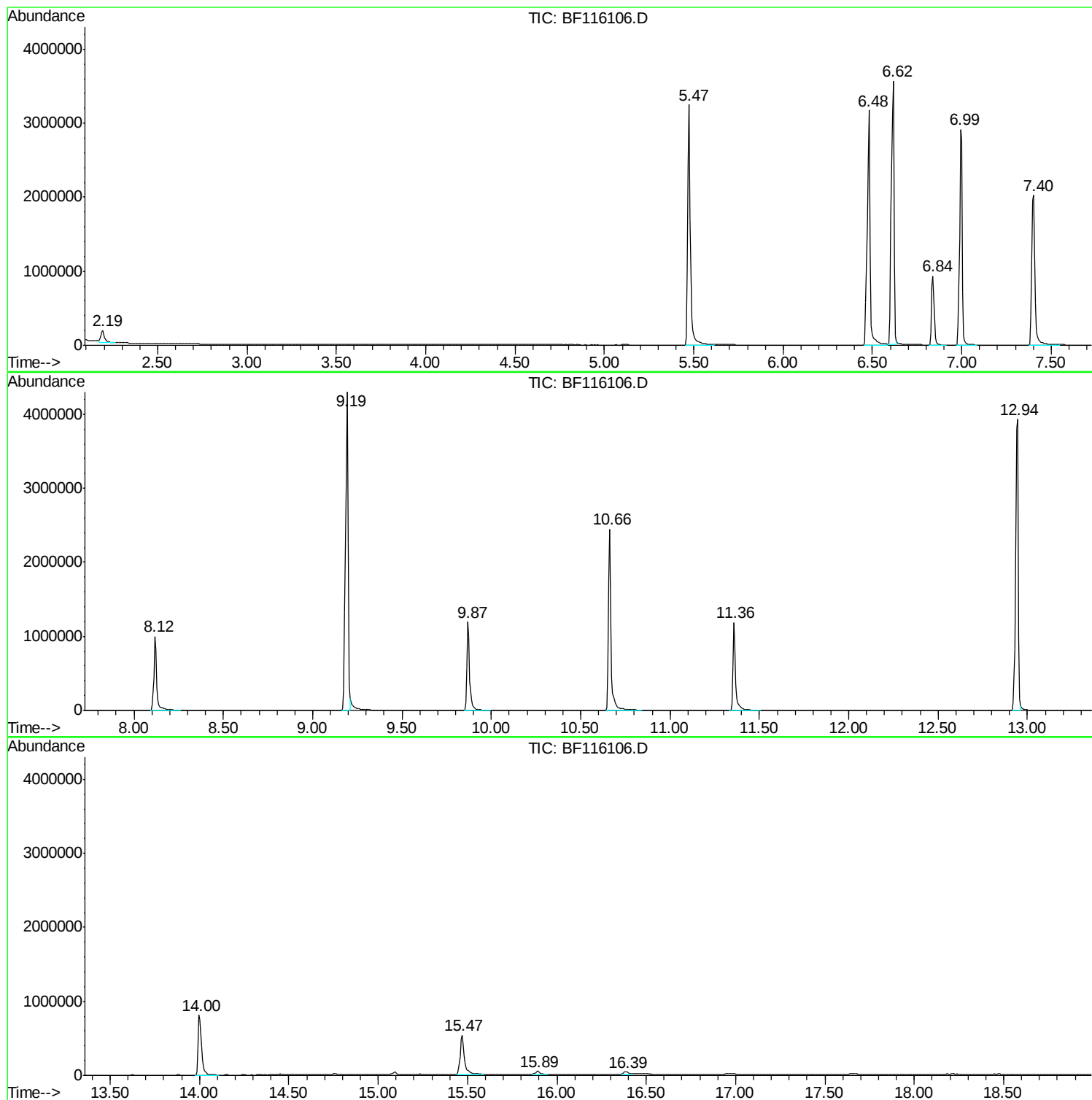
Sum of corrected areas: 33882116

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.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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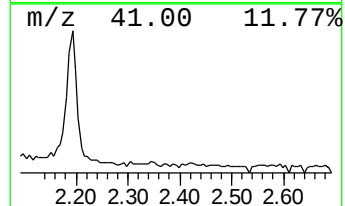
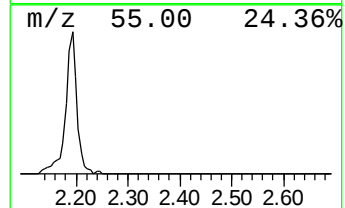
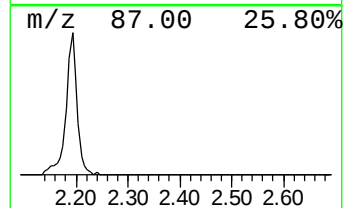
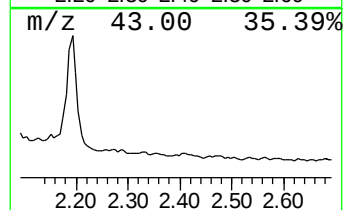
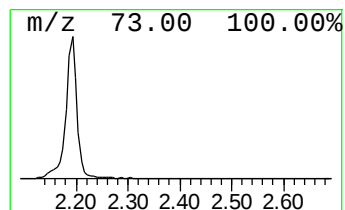
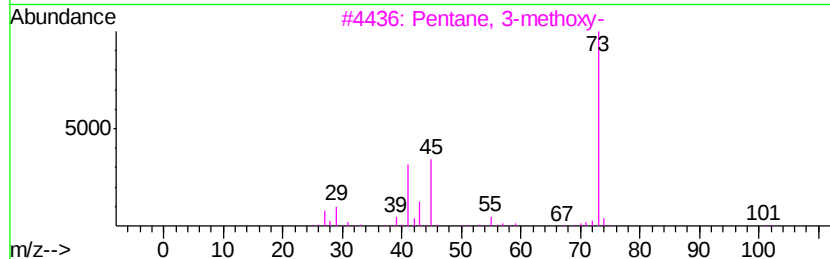
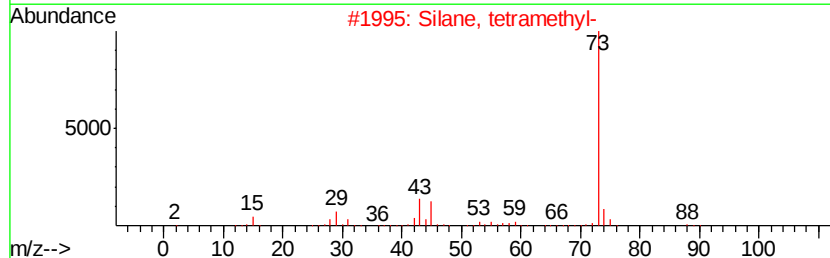
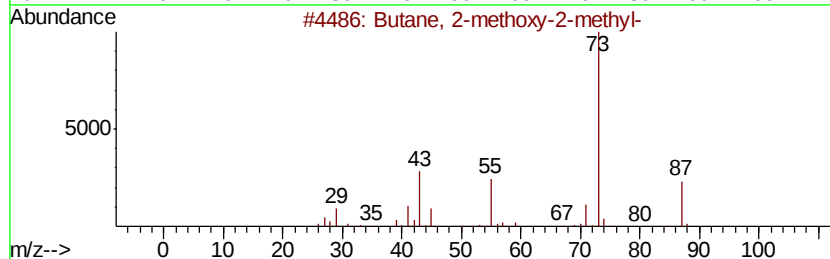
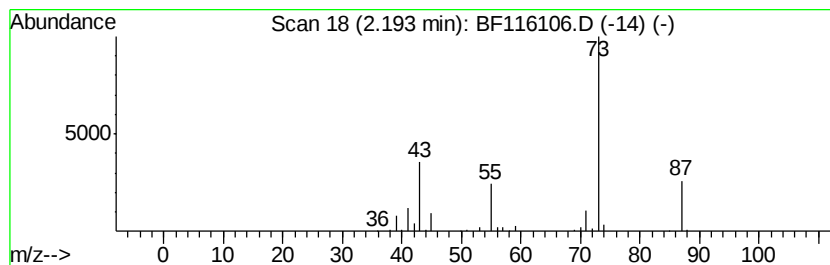
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	5.72 ng	243492	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	59
2			Silane, tetramethyl-	88	C4H12Si	000075-76-3	43
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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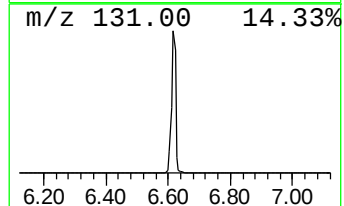
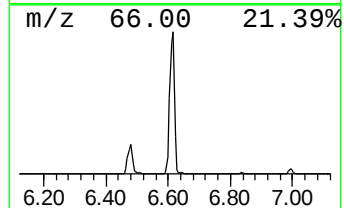
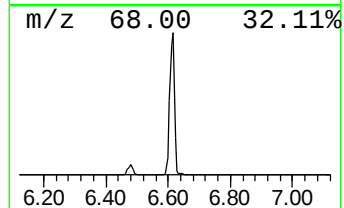
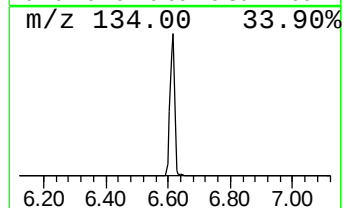
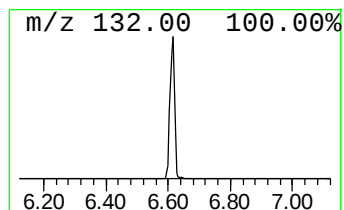
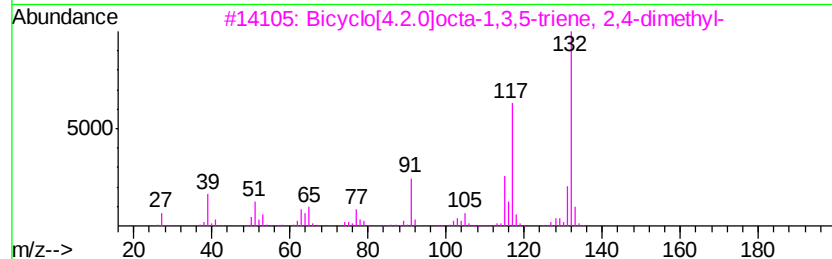
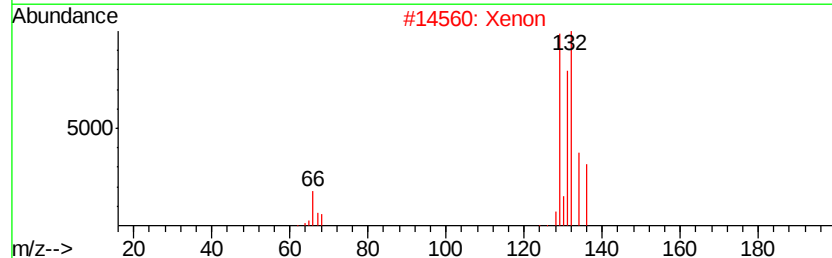
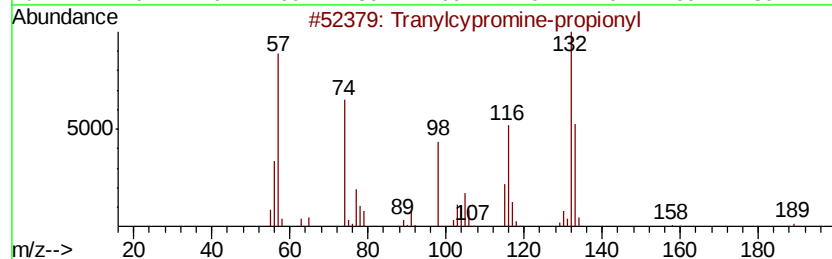
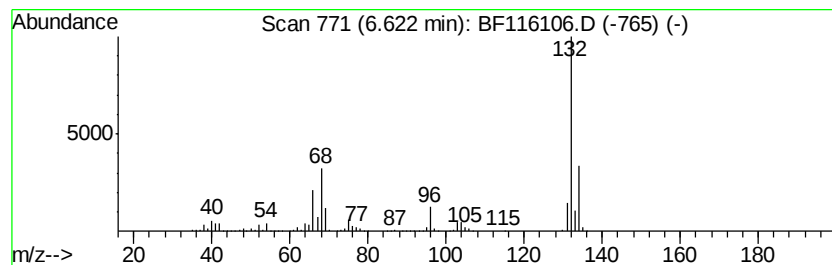
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Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	83.04 ng	3536470	1,4-Dichlorobenzene-d4	6.84

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	12
2		Xenon	132	Xe	007440-63-3	9
3		Bicvclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4		Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.19	5.7	ng	243492	1	6.84	851768	20.0
unknown6.62	6.62	83.0	ng	3536470	1	6.84	851768	20.0