

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121817\
 Data File : BF101509.D
 Acq On : 19 Dec 2017 5:38
 Operator : SJ/JU
 Sample : I6910-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121317-DBRI-E-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:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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	5.022	496	499	510	rBV	55614	81953	2.62%	0.416%
2	5.434	566	569	585	rBV	815506	1053706	33.64%	5.353%
3	6.457	740	743	761	rBV2	488436	733510	23.42%	3.726%
4	6.581	761	764	783	rVB	1870463	1904923	60.82%	9.677%
5	6.810	799	803	811	rBV	845738	756866	24.17%	3.845%
6	6.963	825	829	846	rBV	2621876	2614221	83.47%	13.280%
7	7.381	895	900	916	rBV	1998174	2021919	64.56%	10.272%
8	8.092	1017	1021	1037	rBV	936340	904249	28.87%	4.594%
9	9.169	1199	1204	1207	rBV	3277319	3132014	100.00%	15.911%
10	9.563	1268	1271	1277	rBV	83346	71700	2.29%	0.364%
11	9.845	1316	1319	1324	rVB	723122	696681	22.24%	3.539%
12	10.645	1451	1455	1468	rBV	996968	1034949	33.04%	5.258%
13	11.339	1570	1573	1582	rBV	698220	683331	21.82%	3.471%
14	11.851	1657	1660	1664	rBV	44817	53551	1.71%	0.272%
15	12.563	1778	1781	1789	rBV2	25233	38897	1.24%	0.198%
16	12.921	1838	1842	1846	rBV	2498590	2339199	74.69%	11.883%
17	13.804	1989	1992	1998	rBV3	51743	56463	1.80%	0.287%
18	13.992	2020	2024	2033	rVB	804360	792606	25.31%	4.027%
19	15.462	2269	2274	2281	rVB	528548	713950	22.80%	3.627%

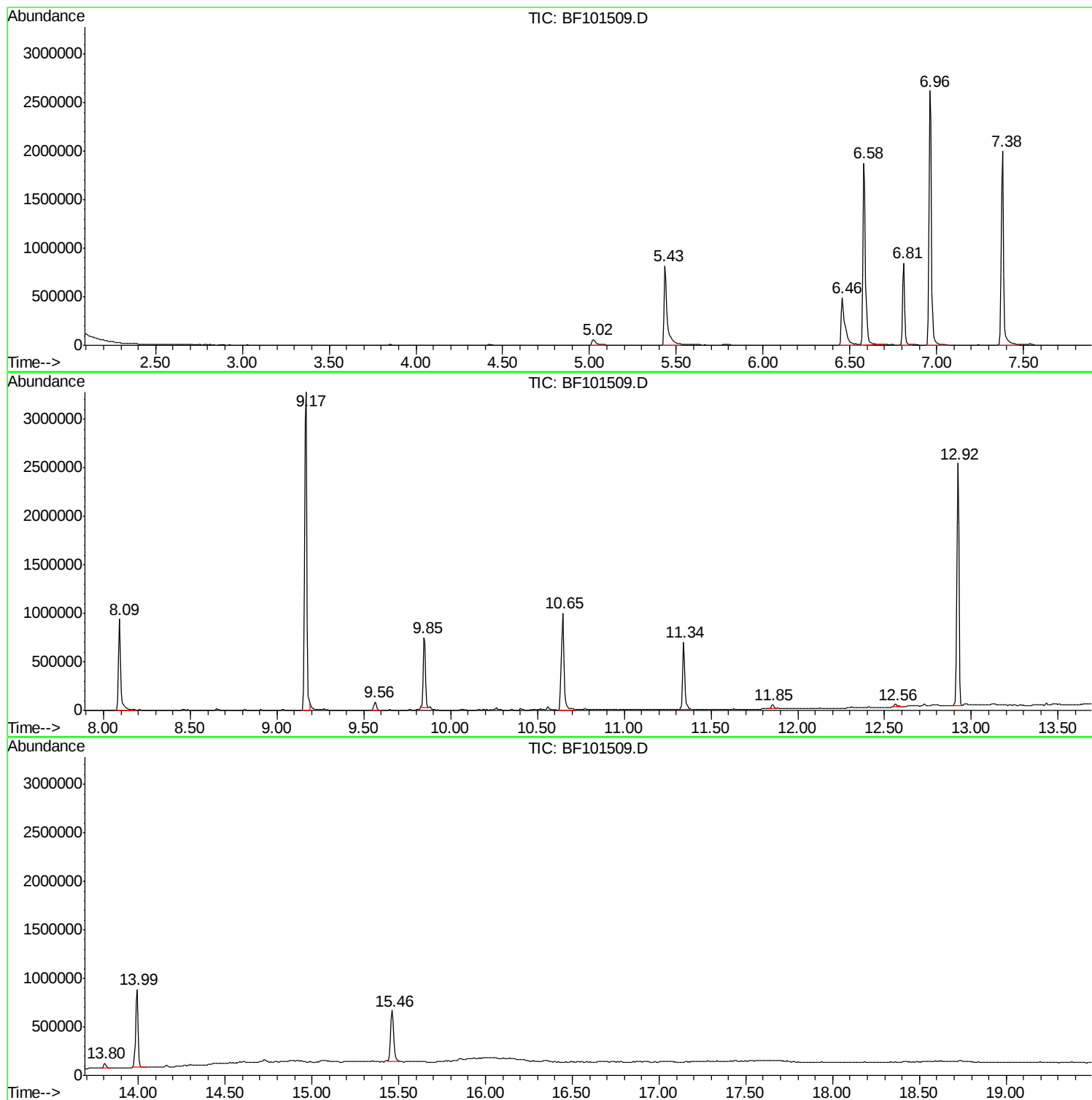
Sum of corrected areas: 19684688

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.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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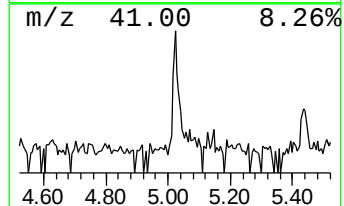
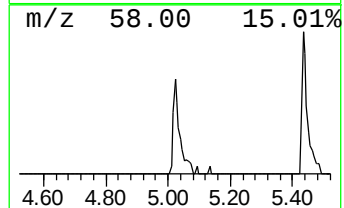
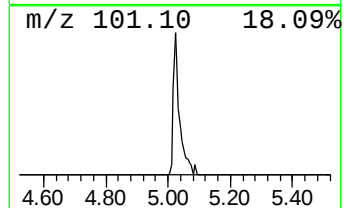
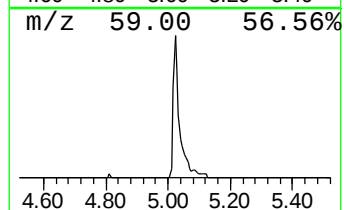
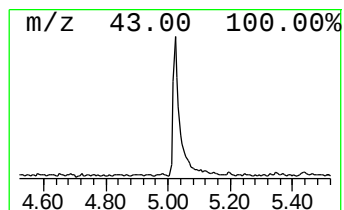
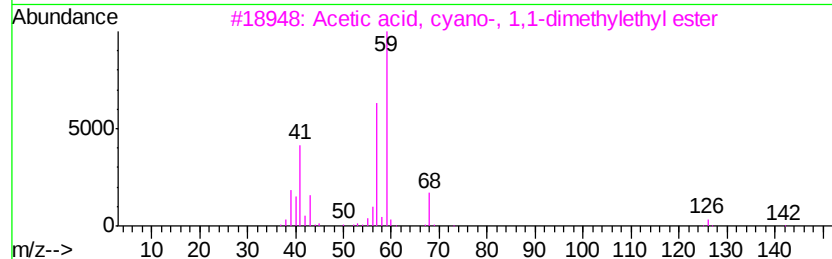
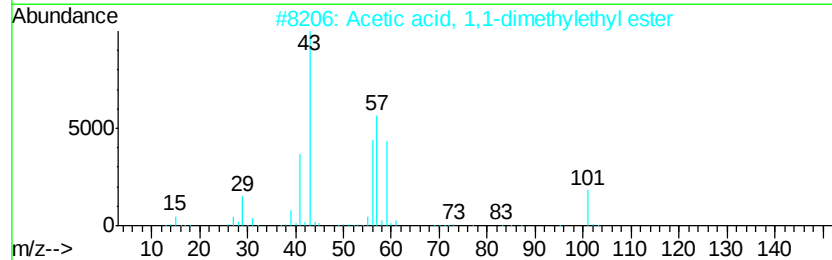
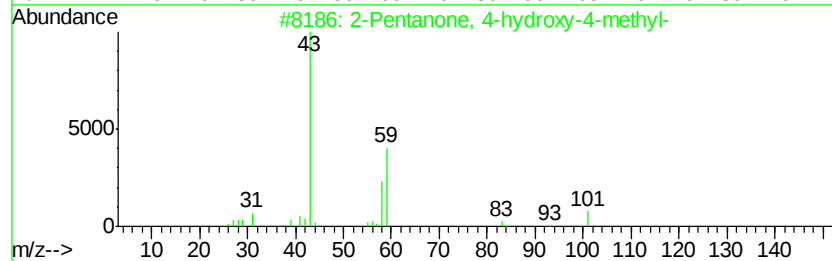
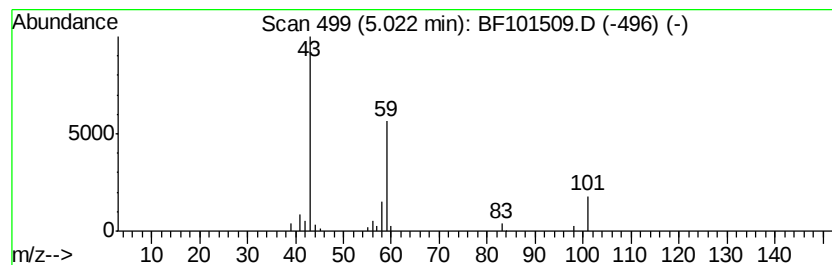
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.02	2.17 ng	81953	1,4-Dichlorobenzene-d4	6.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			5-Hexen-2-one	98	C6H10O	000109-49-9	9
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9



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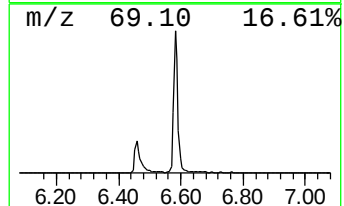
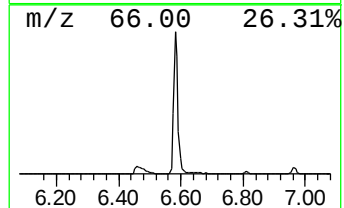
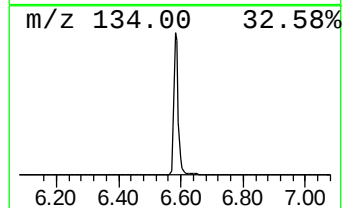
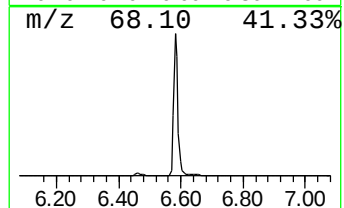
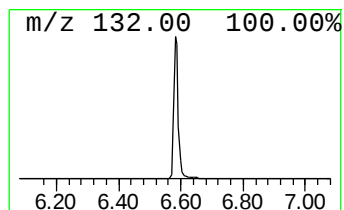
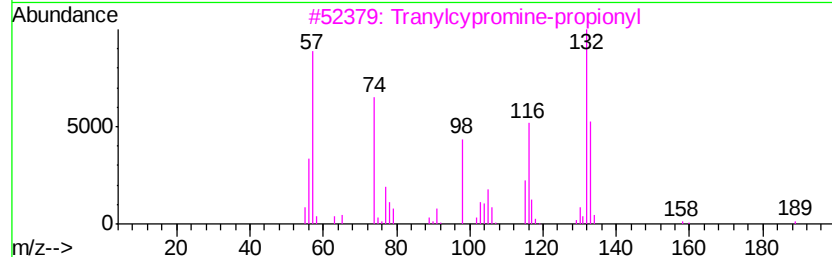
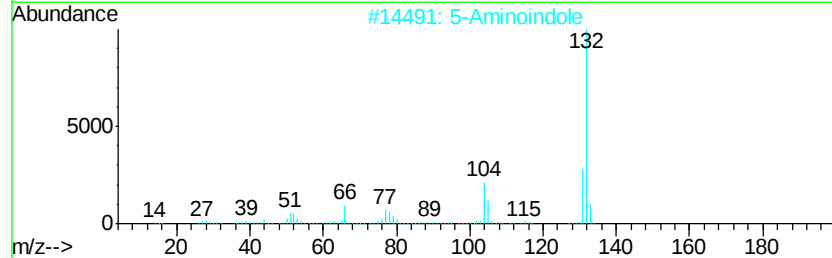
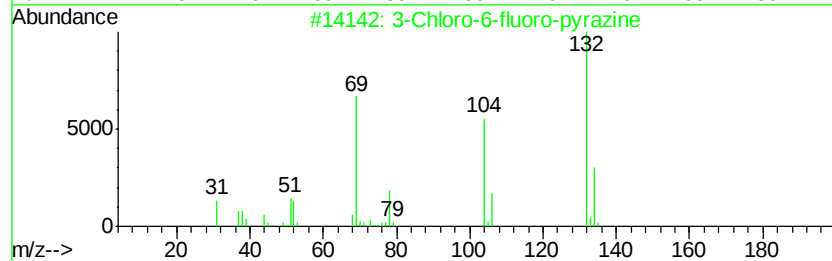
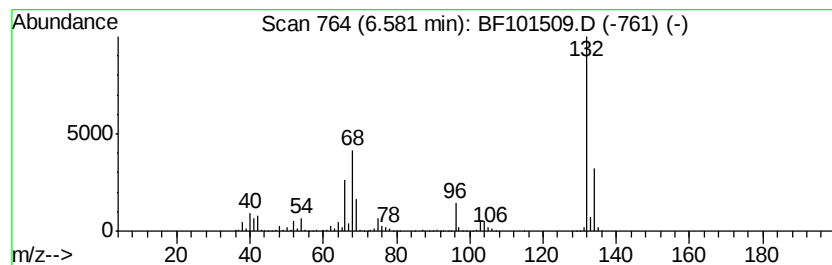
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Peak Number 2 unknown6.58 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	50.34 ng	1904920	1,4-Dichlorobenzene-d4	6.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3	Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5	Aminoindole	132	C8H8N2	005192-03-0	12
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.02	2.2	ng	81953	1	6.81	756866	20.0
unknown6.58	6.58	50.3	ng	1904920	1	6.81	756866	20.0