

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060917\
 Data File : BF095826.D
 Acq On : 9 Jun 2017 17:44
 Operator : SJ/MA
 Sample : PB99623BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99623BL

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-BF060517.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	4.592	506	511	539	rBV2	84115	246426	11.46%	1.539%
2	5.275	622	627	653	rBV	978674	1725244	80.21%	10.773%
3	6.928	903	908	920	rBV	1074839	1635701	76.04%	10.213%
4	7.022	920	924	935	rVB	1304095	1880856	87.44%	11.744%
5	7.369	978	983	994	rBV	251457	343975	15.99%	2.148%
6	7.604	1018	1023	1047	rBV	1037229	1406472	65.39%	8.782%
7	8.280	1133	1138	1161	rBV	690624	1039896	48.34%	6.493%
8	9.398	1323	1328	1343	rBV	338890	479792	22.31%	2.996%
9	11.180	1623	1631	1645	rBV	1441444	2150998	100.00%	13.431%
10	12.221	1802	1808	1819	rBV2	371688	535053	24.87%	3.341%
11	13.515	2023	2028	2047	rBV	750739	1069180	49.71%	6.676%
12	14.615	2210	2215	2228	rBV	398375	541470	25.17%	3.381%
13	16.992	2613	2619	2623	rBV	1708016	2059647	95.75%	12.861%
14	18.333	2842	2847	2857	rBV	412468	527579	24.53%	3.294%
15	19.997	3126	3130	3142	rBV	253066	372963	17.34%	2.329%

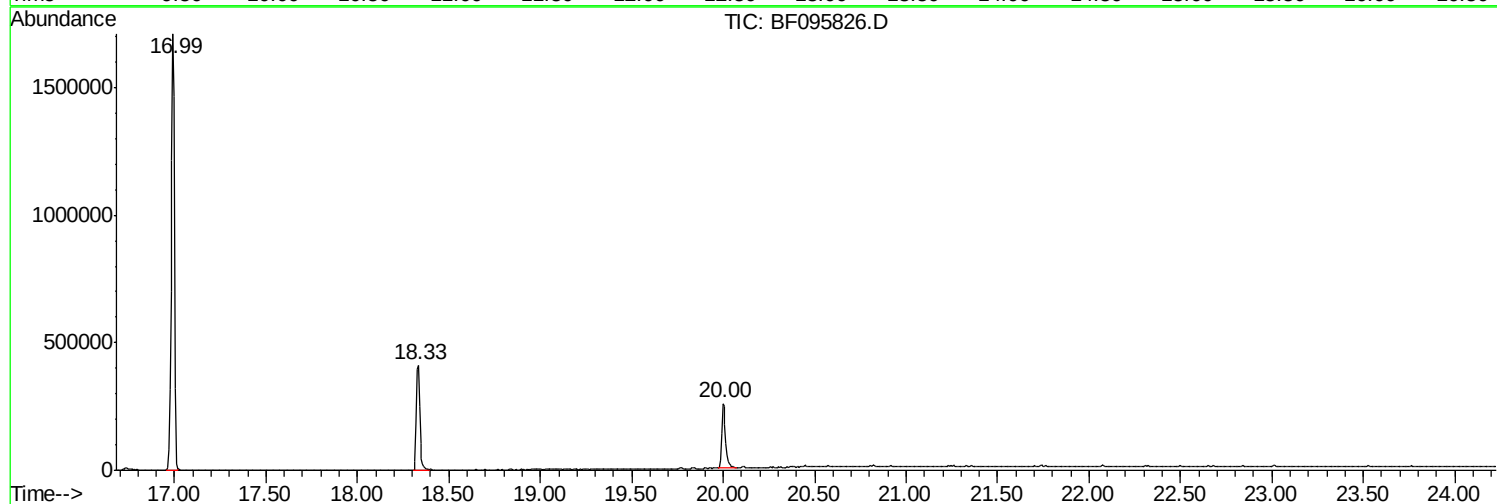
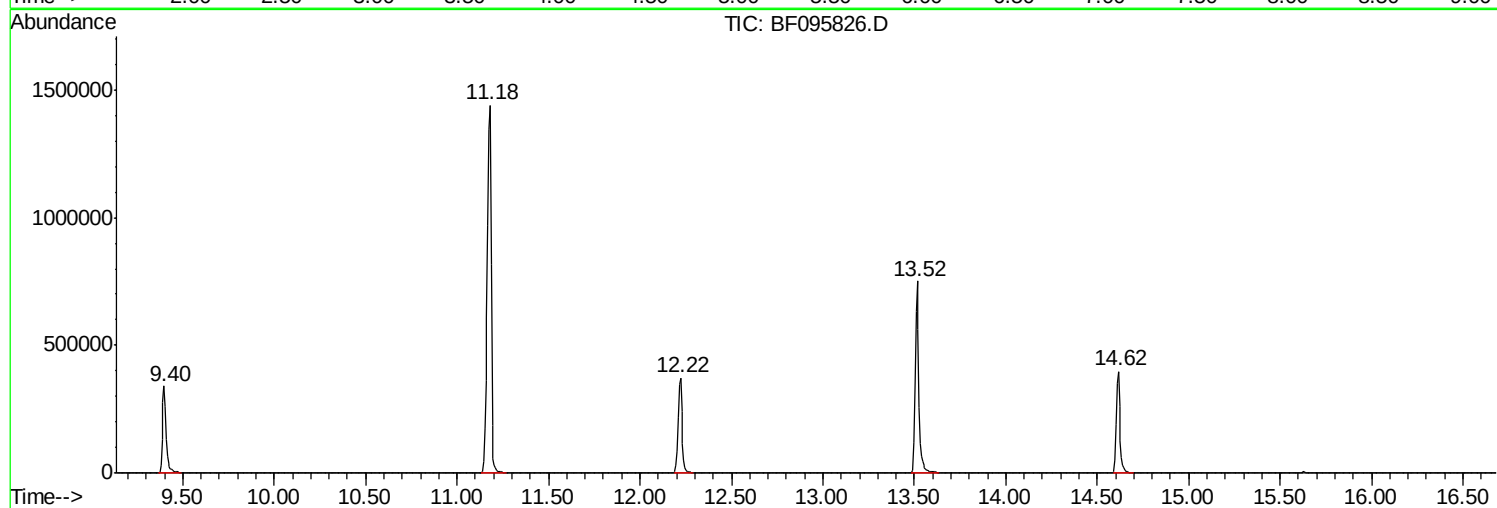
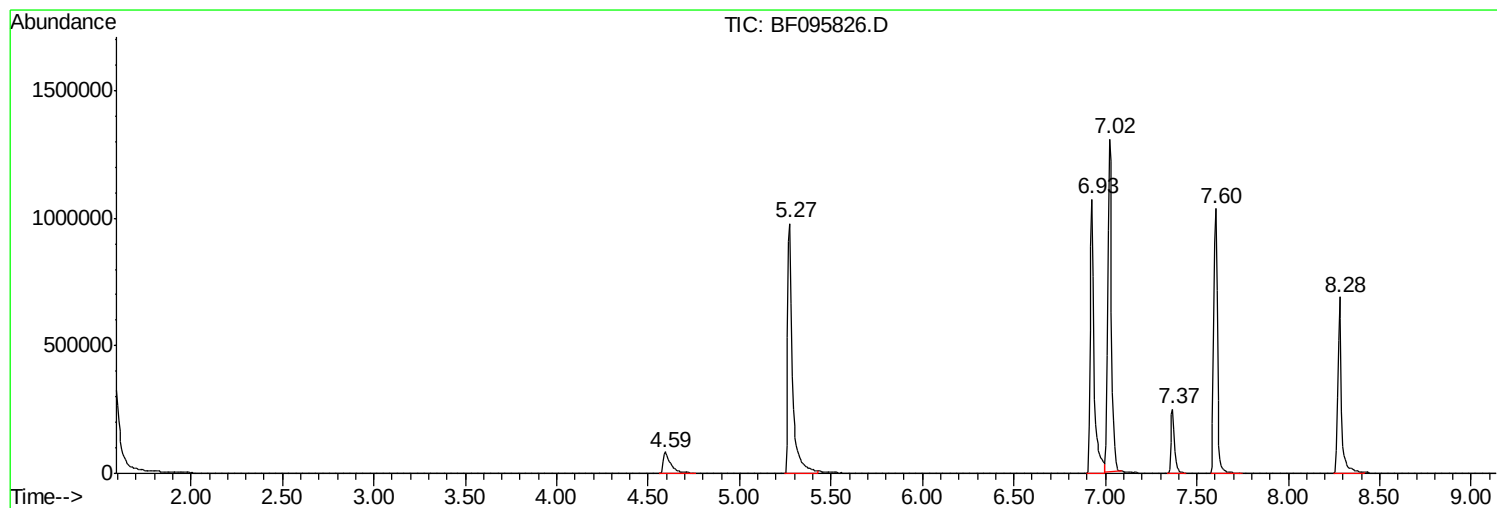
Sum of corrected areas: 16015252

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.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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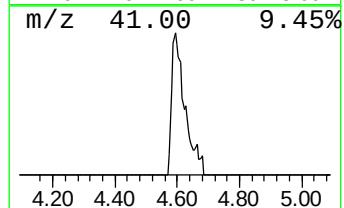
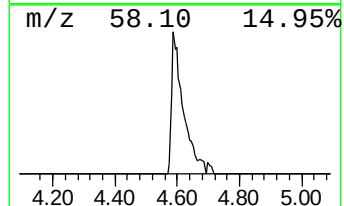
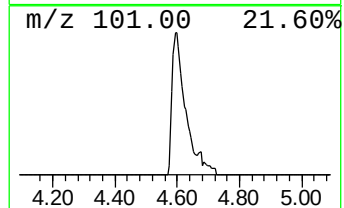
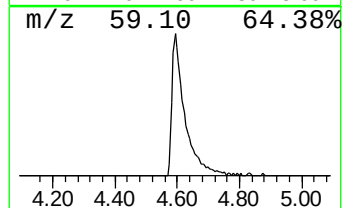
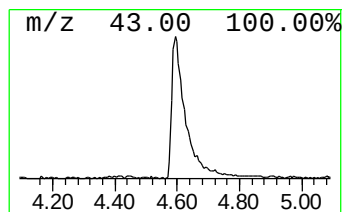
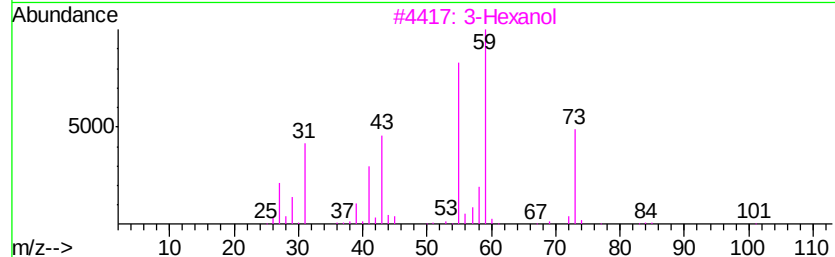
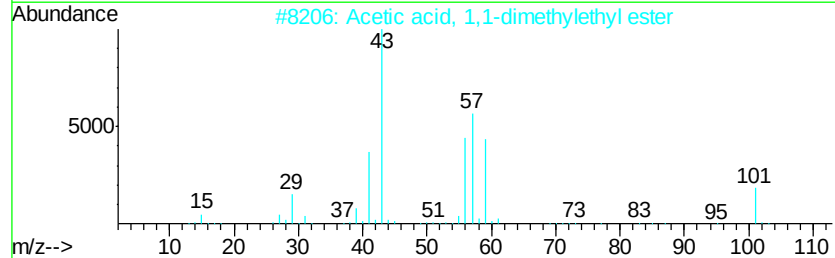
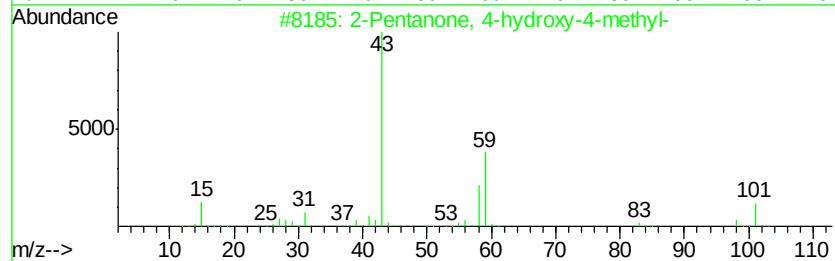
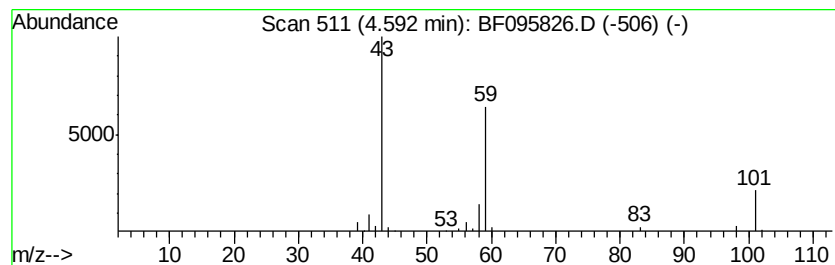
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.59	14.33 ng	246426	1,4-Dichlorobenzene-d4	7.37

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			3-Hexanol	102	C6H14O	000623-37-0	25
4			1,2-Butanediol	90	C4H10O2	000584-03-2	23
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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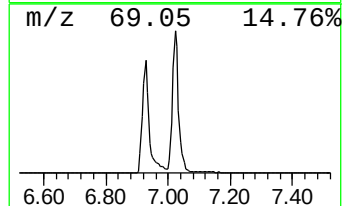
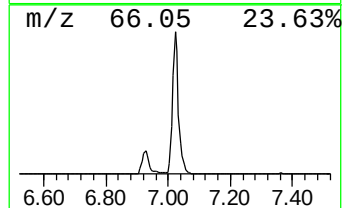
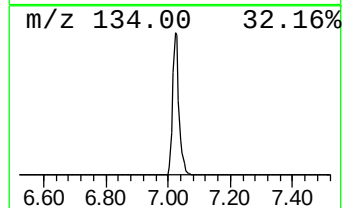
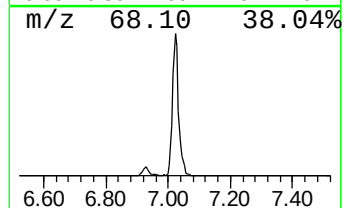
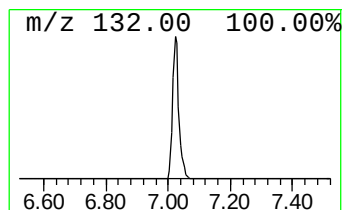
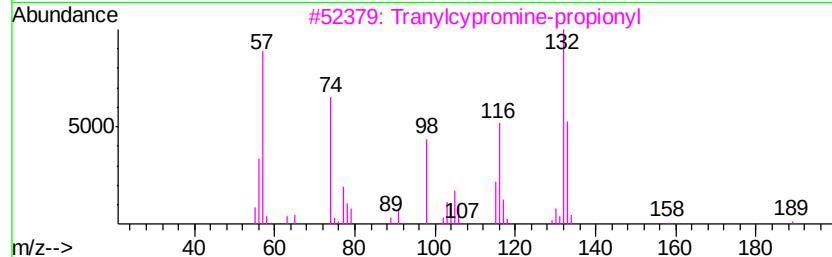
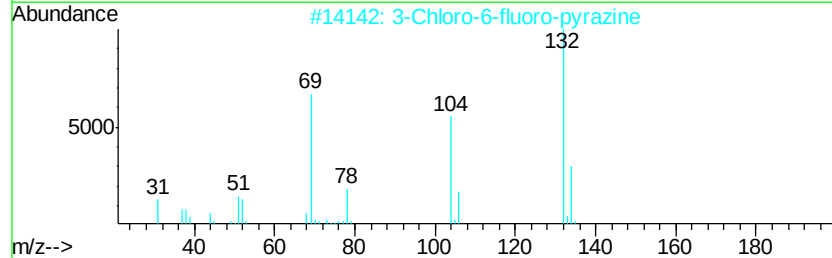
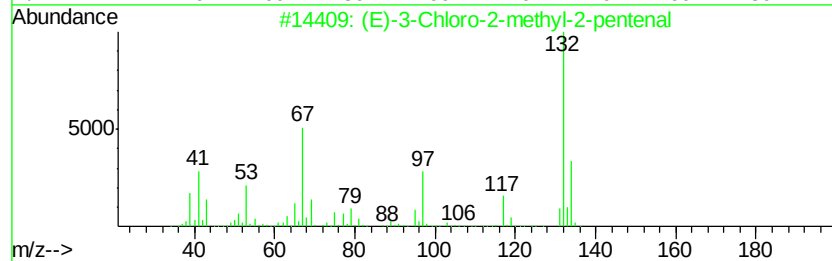
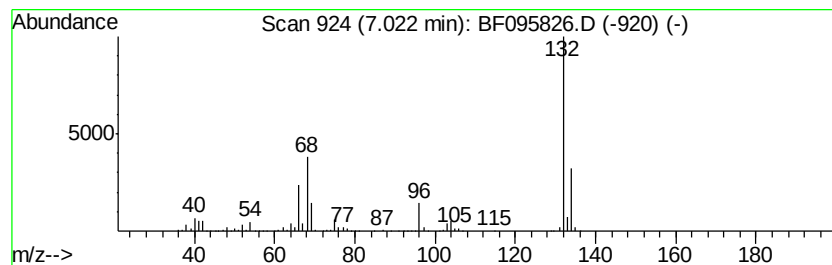
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Peak Number 2 unknown7.02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.02	109.36 ng	1880860	1,4-Dichlorobenzene-d4	7.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
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
2-Pentanone, 4-hy...	4.59	14.3	ng	246426	1	7.37	343975	20.0
unknown7.02	7.02	109.4	ng	1880860	1	7.37	343975	20.0