

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051517\
 Data File : BF095126.D
 Acq On : 15 May 2017 15:28
 Operator : SJ/MA
 Sample : PB98943BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB98943BL

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-BF050217.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	2.013	9	13	25	rVB2	15128	32027	1.07%	0.151%
2	5.160	544	548	569	rBV	244073	559171	18.60%	2.643%
3	5.760	646	650	682	rBV	1176762	2296343	76.40%	10.852%
4	7.325	911	916	933	rBV	1181358	1912011	63.62%	9.036%
5	7.448	933	937	955	rVB	1668970	2324482	77.34%	10.985%
6	7.783	991	994	1008	rVB	354983	462070	15.37%	2.184%
7	8.019	1030	1034	1053	rBV	1481584	1796909	59.79%	8.492%
8	8.677	1142	1146	1165	rBV	874455	1289931	42.92%	6.096%
9	9.801	1332	1337	1357	rBV	412705	679244	22.60%	3.210%
10	11.565	1632	1637	1663	rBV	2263350	3005493	100.00%	14.204%
11	12.624	1813	1817	1839	rBV2	406573	686654	22.85%	3.245%
12	13.924	2033	2038	2064	rBV	932213	1576406	52.45%	7.450%
13	15.036	2222	2227	2241	rBV2	423769	726547	24.17%	3.434%
14	17.353	2615	2621	2634	rBV	2202073	2701972	89.90%	12.769%
15	18.700	2846	2850	2871	rBV	371343	620602	20.65%	2.933%
16	20.377	3131	3135	3146	rBV	274176	490261	16.31%	2.317%

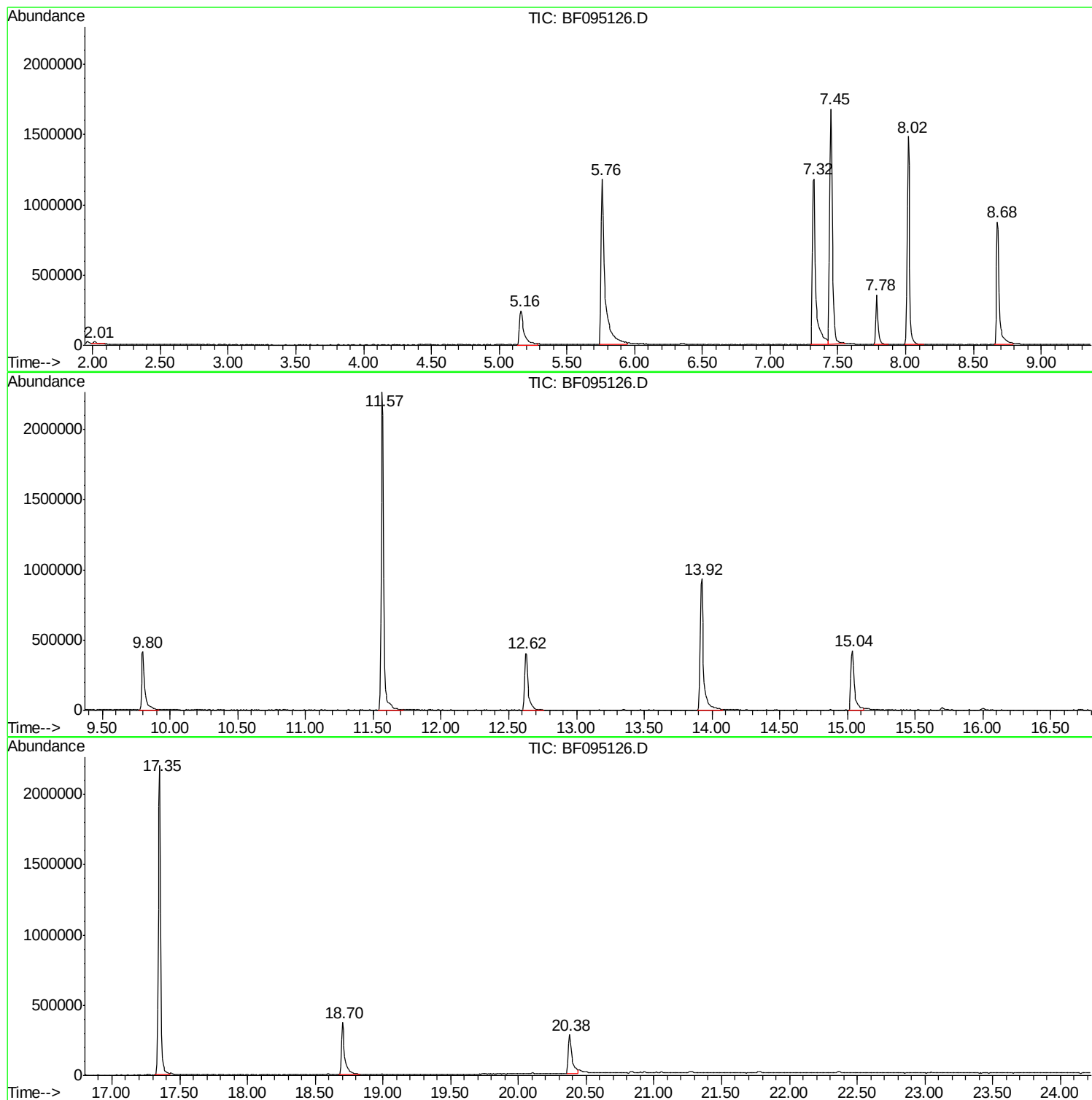
Sum of corrected areas: 21160123

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.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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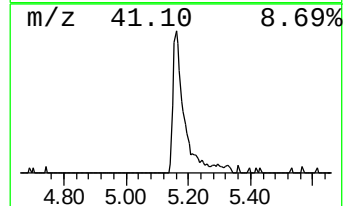
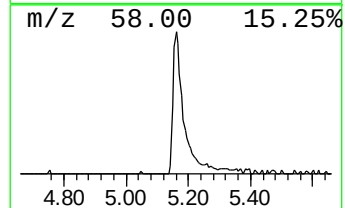
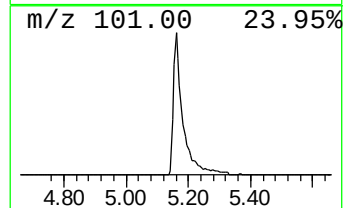
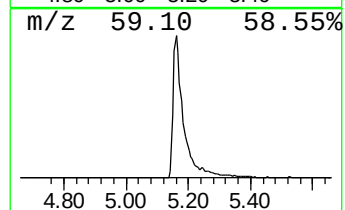
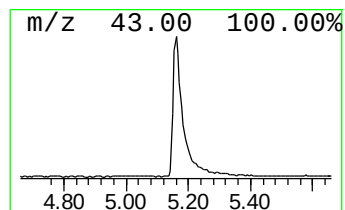
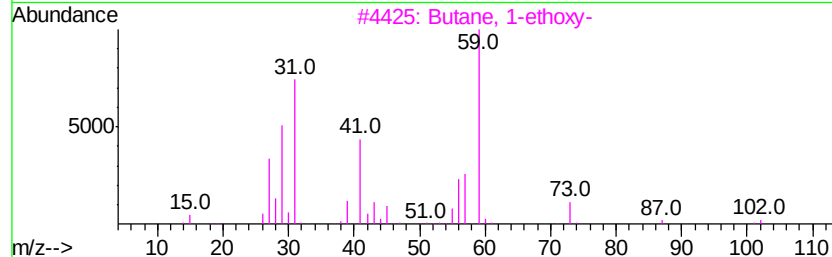
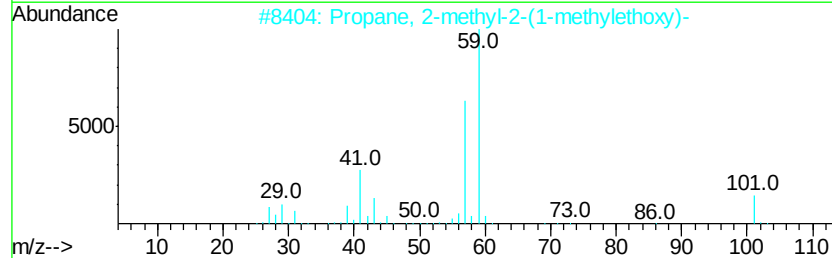
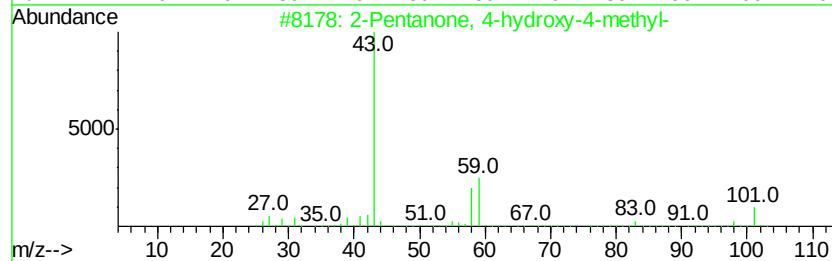
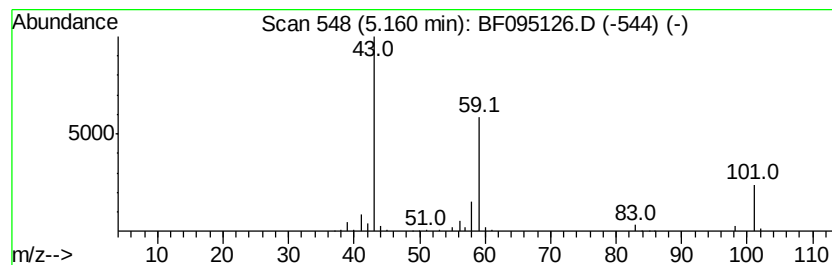
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	24.20 ng	559171	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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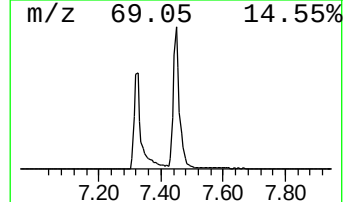
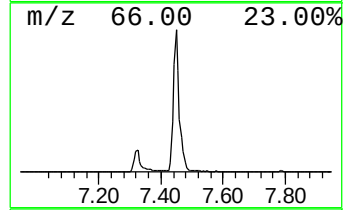
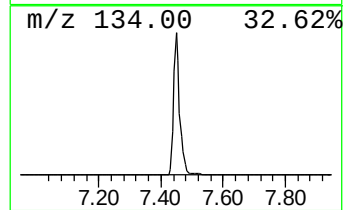
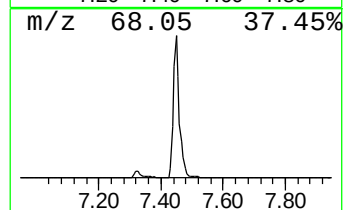
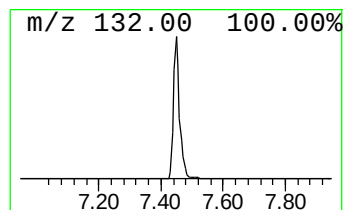
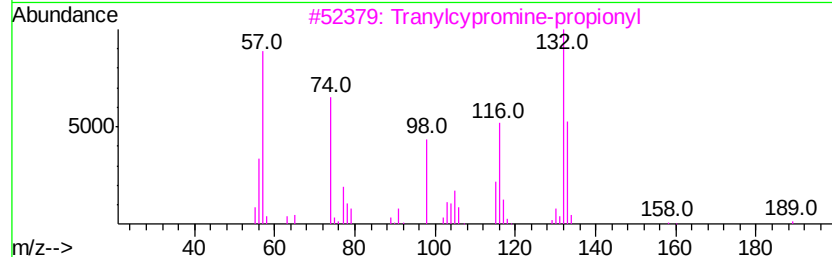
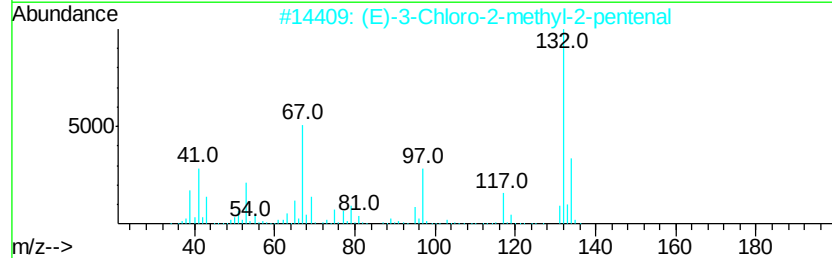
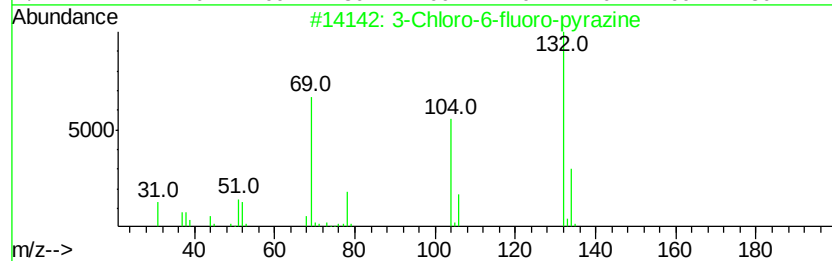
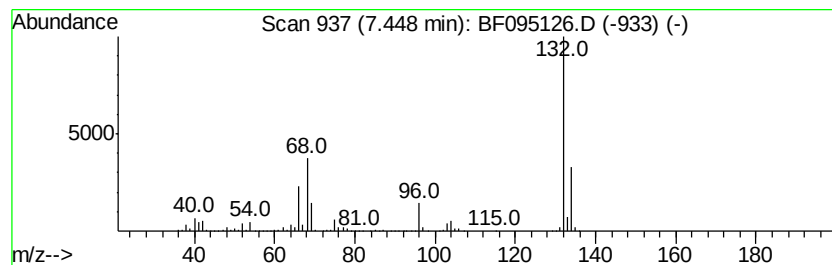
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Peak Number 2 unknown7.45 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.45	100.61 ng	2324480	1,4-Dichlorobenzene-d4	7.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.16	24.2	ng	559171	1	7.78	462070	20.0
unknown7.45	7.45	100.6	ng	2324480	1	7.78	462070	20.0