

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122017\
 Data File : BF101591.D
 Acq On : 21 Dec 2017 6:06
 Operator : SJ/JU
 Sample : I6971-02
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121517-TIDO-F-D-S

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	5.010	494	497	511	rBV	78122	125951	3.66%	0.548%
2	5.416	563	566	593	rBV	1102845	1523655	44.26%	6.628%
3	6.434	736	739	758	rBV	617411	965241	28.04%	4.199%
4	6.569	758	762	769	rBV	2131106	2250149	65.37%	9.789%
5	6.793	797	800	810	rBV	973271	905147	26.29%	3.938%
6	6.951	823	827	843	rBV	2920271	2845950	82.67%	12.381%
7	7.363	893	897	918	rBV	2230539	2462600	71.54%	10.713%
8	8.075	1015	1018	1036	rBV	1047585	1078049	31.32%	4.690%
9	9.151	1196	1201	1204	rBV	3685605	3442437	100.00%	14.976%
10	9.834	1313	1317	1330	rVB	941150	963653	27.99%	4.192%
11	10.628	1448	1452	1463	rBV	1064926	1164848	33.84%	5.068%
12	11.322	1566	1570	1580	rBV	864306	819578	23.81%	3.565%
13	12.710	1802	1806	1809	rBV	48627	40998	1.19%	0.178%
14	12.904	1834	1839	1842	rBV	2738349	2495365	72.49%	10.856%
15	13.792	1985	1990	2001	rBV3	25677	60931	1.77%	0.265%
16	13.969	2016	2020	2031	rBV	932684	889857	25.85%	3.871%
17	15.433	2265	2269	2283	rVB	676224	915743	26.60%	3.984%
18	16.304	2412	2417	2422	rBV5	19731	36400	1.06%	0.158%

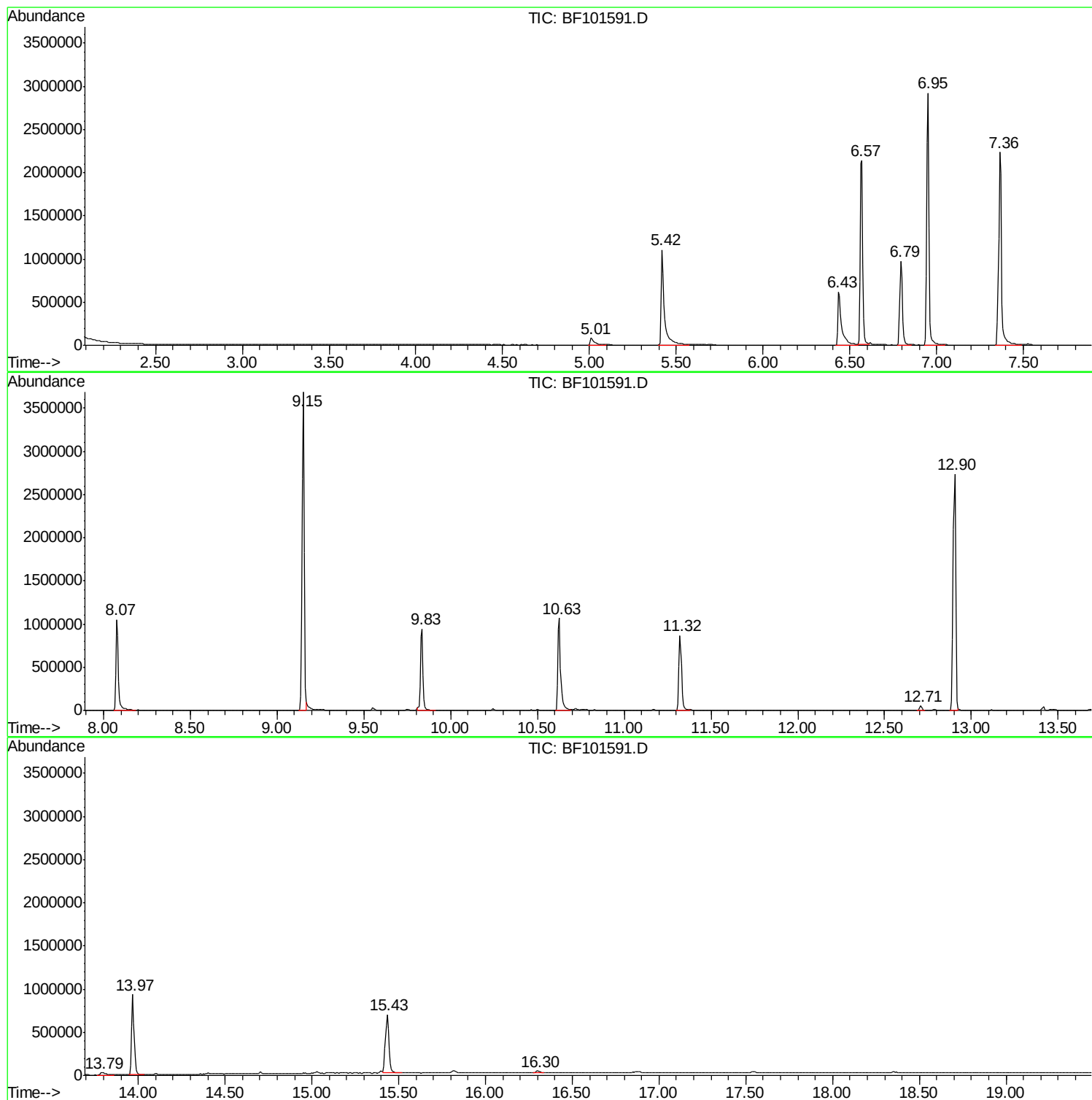
Sum of corrected areas: 22986552

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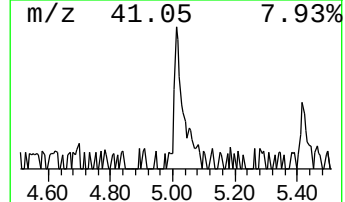
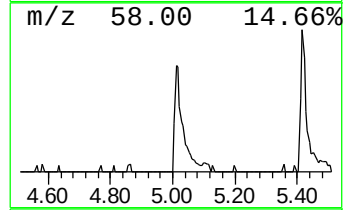
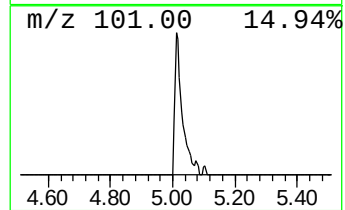
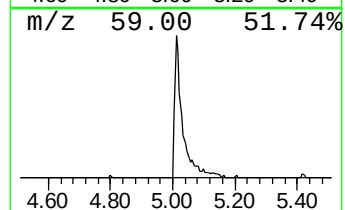
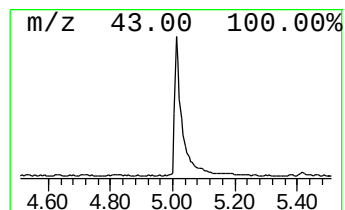
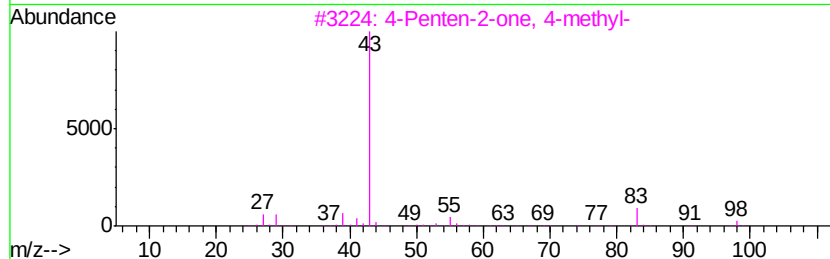
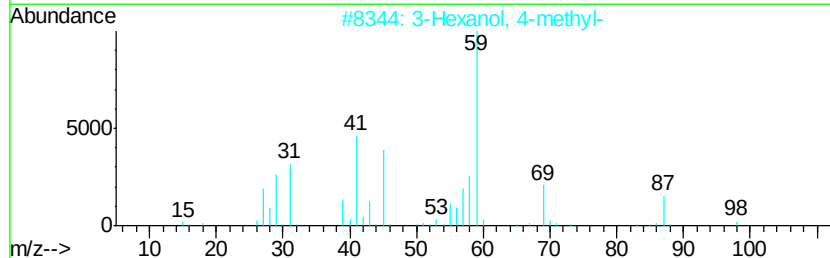
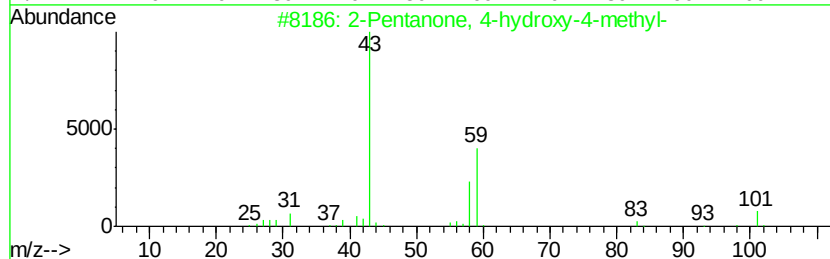
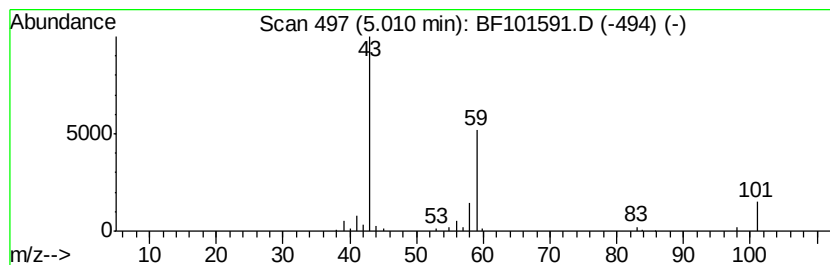
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.01	2.78 ng	125951	1,4-Dichlorobenzene-d4	6.79

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	4-Penten-2-one, 4-methyl-	98	C6H10O	003744-02-3	9		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	Diacetamide	101	C4H7NO2	000625-77-4	9		



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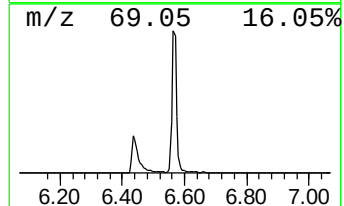
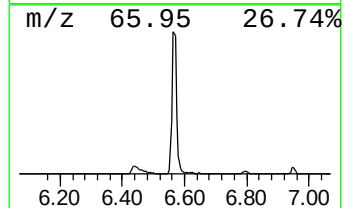
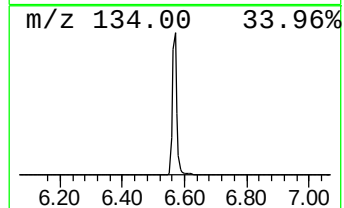
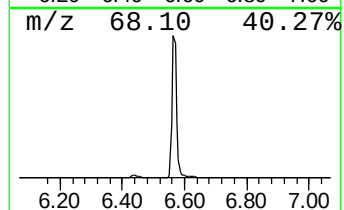
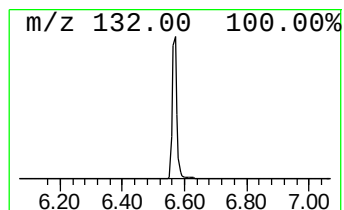
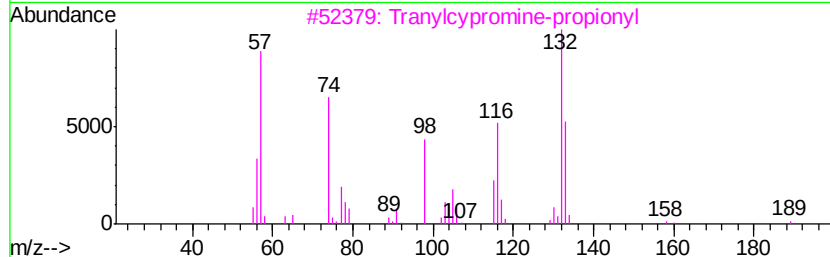
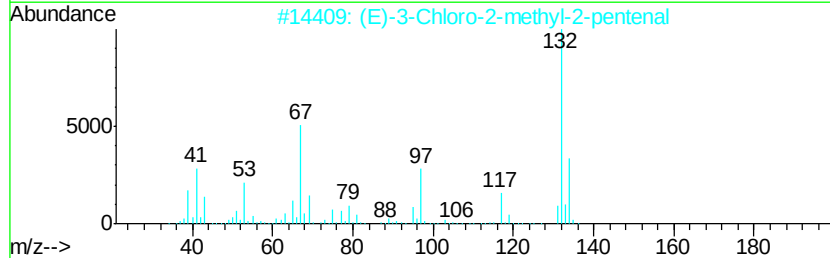
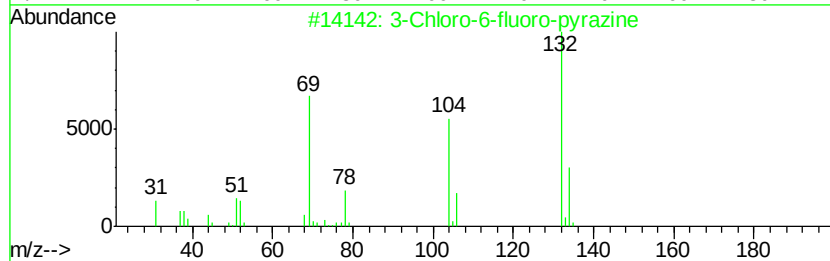
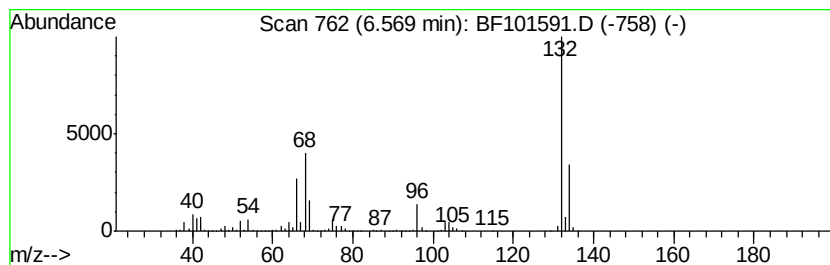
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Peak Number 2 unknown6.57 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.57	49.72 ng	2250150	1,4-Dichlorobenzene-d4	6.79

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
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		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.01	2.8	ng	125951	1	6.79	905147	20.0
unknown6.57	6.57	49.7	ng	2250150	1	6.79	905147	20.0