

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF081015\
 Data File : BF080955.D
 Acq On : 10 Aug 2015 17:04
 Operator : TP/UM
 Sample : PB84929BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB84929BL

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0
 Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080715.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	4.304	191	194	198	rBV	44836	53709	2.23%	0.264%
2	4.979	248	253	265	rBV	804196	1540826	63.89%	7.565%
3	5.390	286	289	291	rBV	1754685	1841130	76.34%	9.039%
4	5.790	322	324	327	rBV	55089	47202	1.96%	0.232%
5	6.419	376	379	382	rBV	1287334	1863053	77.25%	9.146%
6	6.556	388	391	393	rBV	1810325	1892861	78.49%	9.293%
7	6.785	409	411	413	rBV	628759	522097	21.65%	2.563%
8	6.945	422	425	427	rBV	1450029	1486940	61.66%	7.300%
9	7.345	457	460	463	rBV	756669	1176198	48.77%	5.774%
10	8.065	521	523	526	rBV	575909	659966	27.37%	3.240%
11	9.150	614	618	620	rBV	2215872	2169383	89.96%	10.650%
12	9.825	674	677	679	rBB	762657	809053	33.55%	3.972%
13	10.613	742	746	748	rBV2	973718	1424909	59.09%	6.995%
14	11.299	803	806	808	rBV	1012185	869786	36.07%	4.270%
15	12.888	942	945	947	rBV	2420518	2411611	100.00%	11.840%
16	13.928	1033	1036	1038	rBV	827502	880819	36.52%	4.324%
17	15.323	1155	1158	1161	rBV	598666	719526	29.84%	3.532%

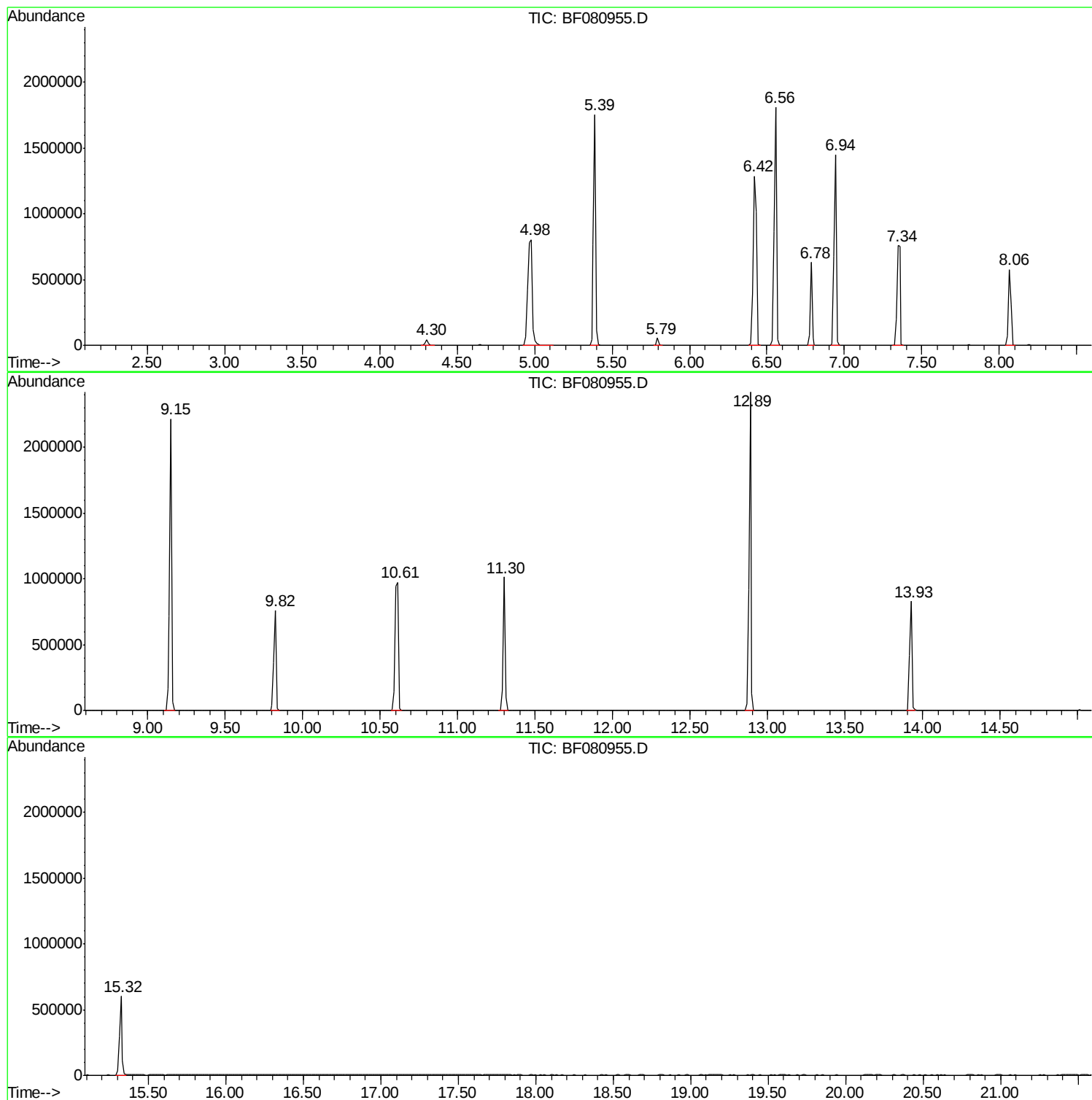
Sum of corrected areas: 20369069

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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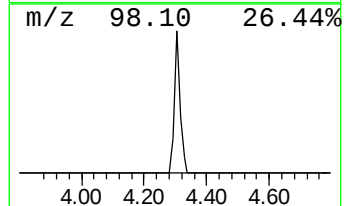
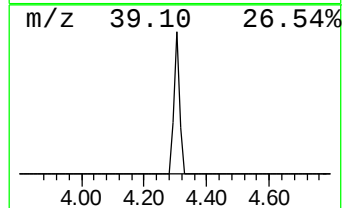
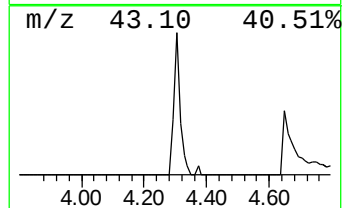
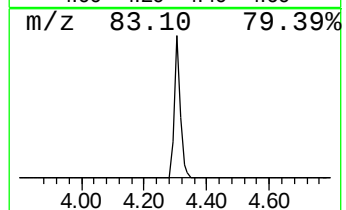
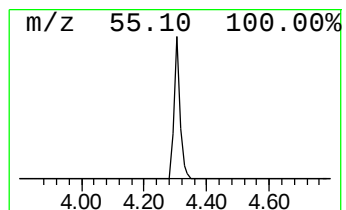
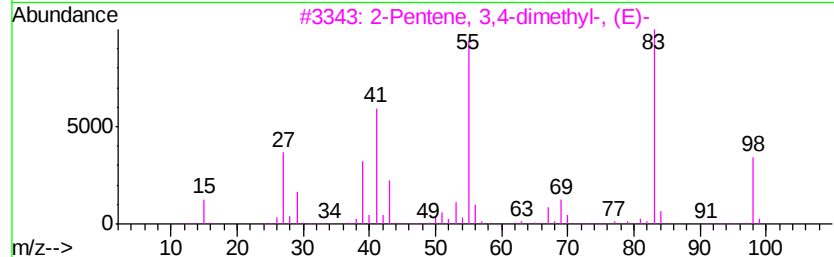
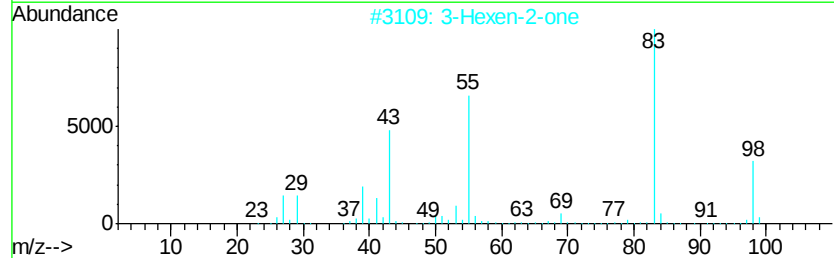
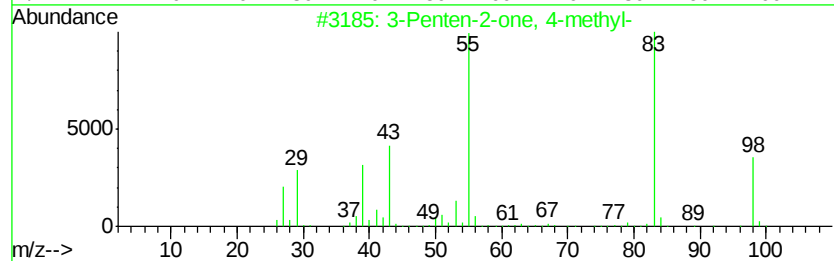
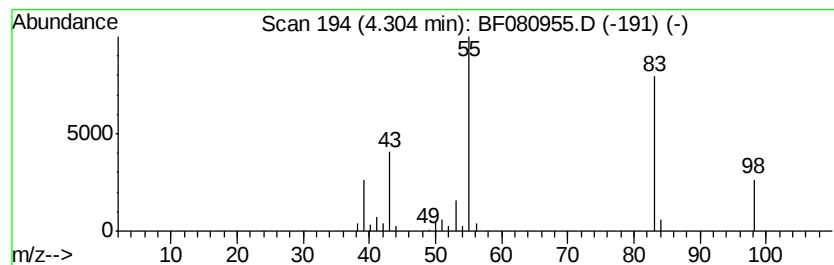
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.30	2.06 ng	53709	1,4-Dichlorobenzene-d4	6.78

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2			3-Hexen-2-one	98	C6H10O	000763-93-9	86
3			2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	86
4			2-Pentene, 2,4-dimethyl-	98	C7H14	000625-65-0	80
5			2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	80



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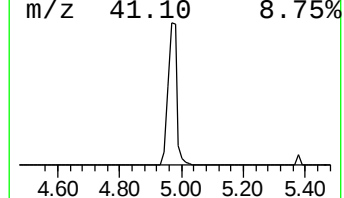
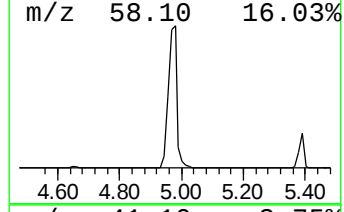
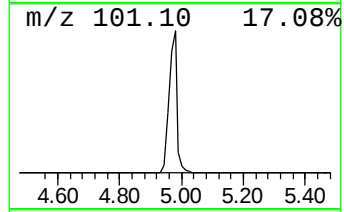
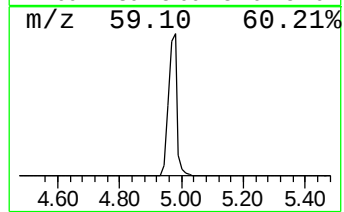
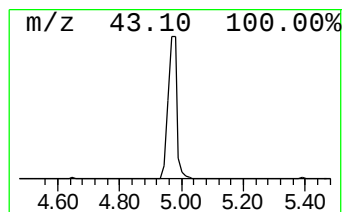
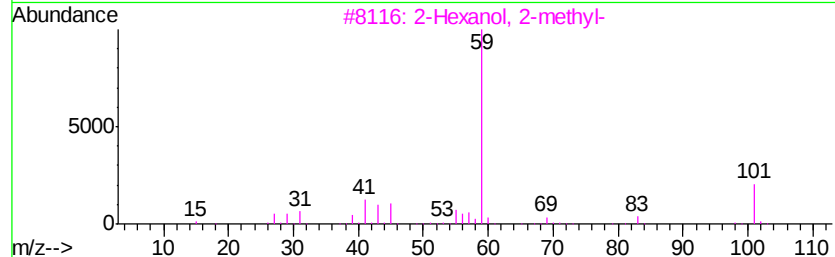
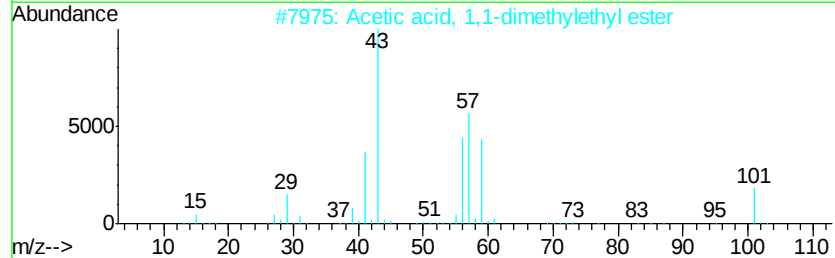
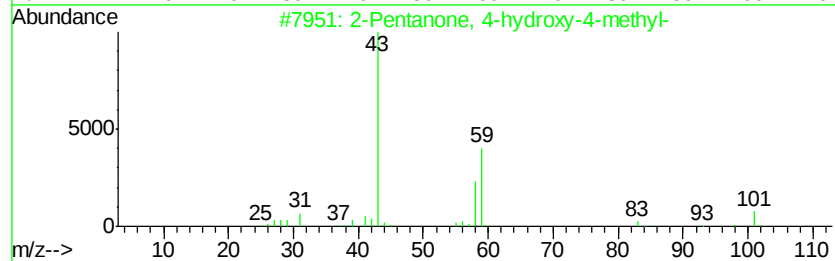
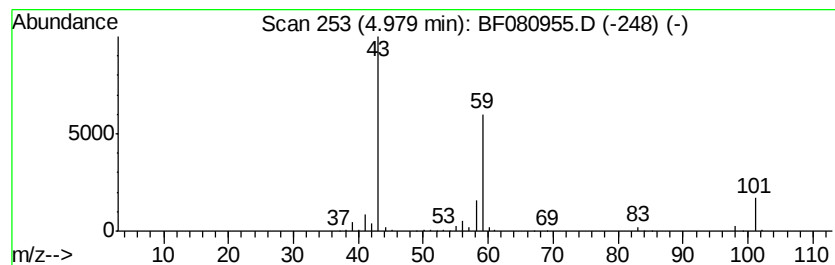
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.98	59.02 ng	1540830	1,4-Dichlorobenzene-d4	6.78

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3		2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
4		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
5		Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
3-Penten-2-one, 4...	4.30	2.1	ng	53709	1	6.78	522097	20.0
2-Pentanone, 4-hy...	4.98	59.0	ng	1540830	1	6.78	522097	20.0