

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF122018\
 Data File : BF111656.D
 Acq On : 20 Dec 2018 23:18
 Operator : JU/SJ
 Sample : J6362-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-4

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF121918.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	2.222	20	23	29	rVB	50366	75079	2.18%	0.245%
2	3.428	225	228	236	rBV	23236	52257	1.52%	0.171%
3	5.116	511	515	520	rBV	129149	160526	4.66%	0.525%
4	5.516	578	583	587	rBV	3008715	3069297	89.07%	10.034%
5	6.522	748	754	758	rBV	3318747	3446016	100.00%	11.265%
6	6.663	774	778	782	rBV	3463251	3219981	93.44%	10.526%
7	6.898	814	818	821	rBV	1499920	1246653	36.18%	4.075%
8	7.051	840	844	848	rBV	2856847	2543122	73.80%	8.314%
9	7.451	907	912	915	rBV	2383395	1993529	57.85%	6.517%
10	8.175	1031	1035	1039	rBV	1818010	1557222	45.19%	5.091%
11	9.251	1213	1218	1221	rBV	4382752	3432584	99.61%	11.222%
12	9.639	1280	1284	1287	rBV	286197	209196	6.07%	0.684%
13	9.933	1330	1334	1337	rVB	1701990	1542646	44.77%	5.043%
14	10.716	1462	1467	1470	rBV	1780800	1658557	48.13%	5.422%
15	11.416	1582	1586	1589	rBV	1357266	1265321	36.72%	4.136%
16	11.933	1669	1674	1679	rBV	91408	103217	3.00%	0.337%
17	12.992	1850	1854	1858	rBV	2670747	2514071	72.96%	8.219%
18	13.892	2004	2007	2012	rBV	118192	134158	3.89%	0.439%
19	14.045	2029	2033	2037	rBV	1572726	1394294	40.46%	4.558%
20	15.521	2279	2284	2288	rBV	809813	971621	28.20%	3.176%

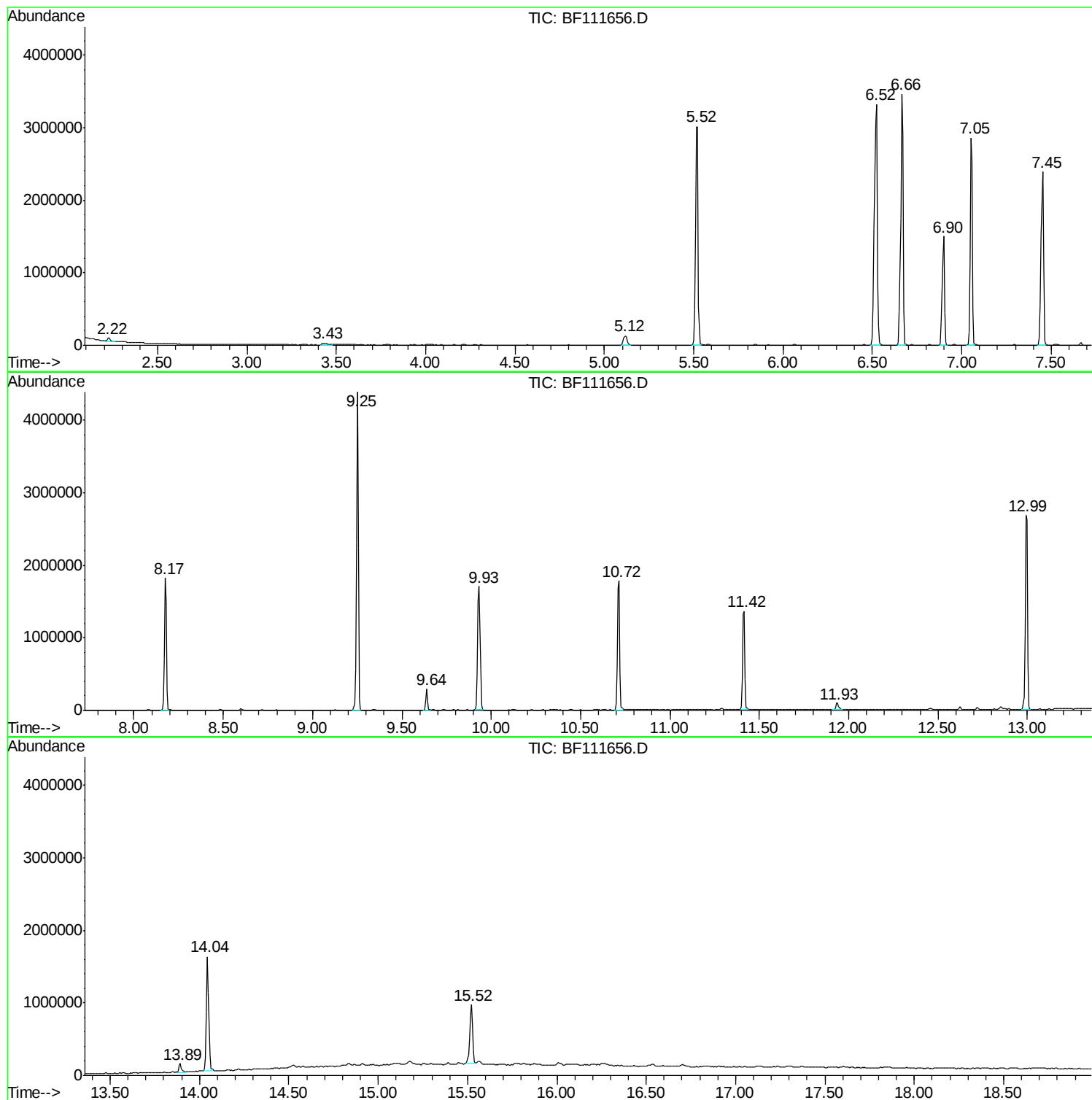
Sum of corrected areas: 30589347

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.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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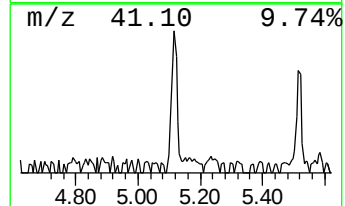
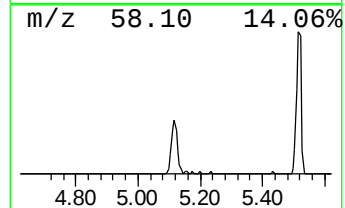
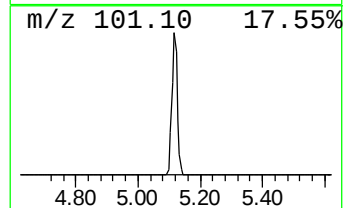
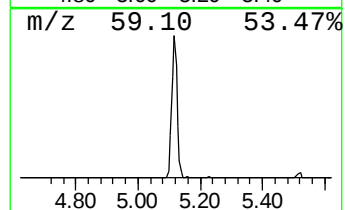
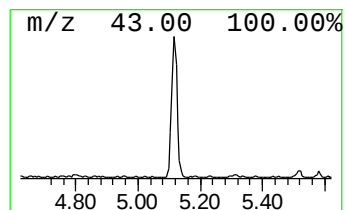
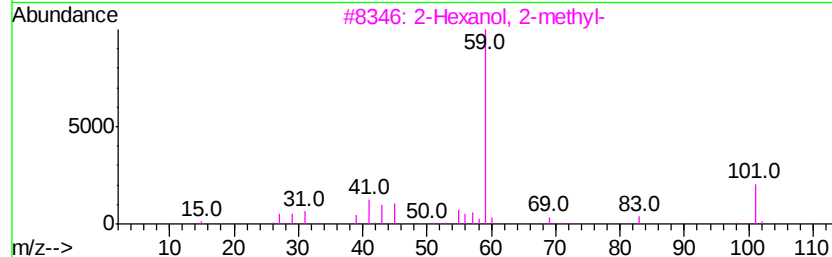
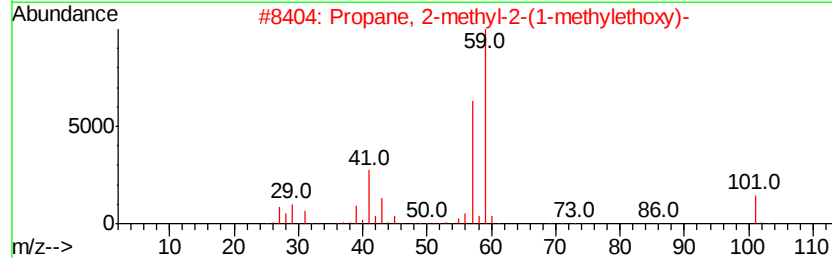
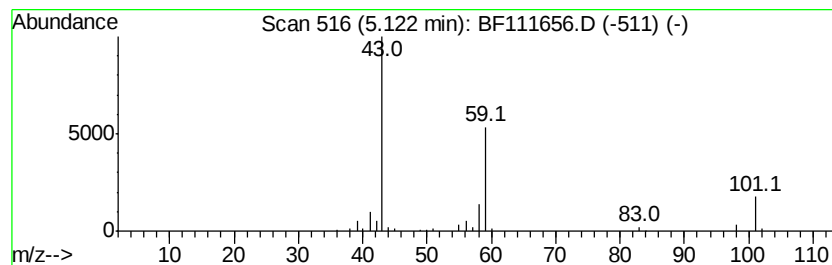
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.12	2.58 ng	160526	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
4			3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	9
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	9



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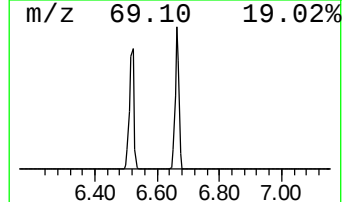
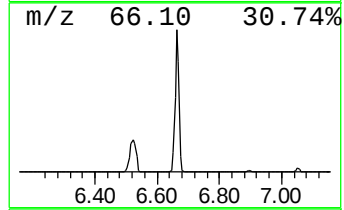
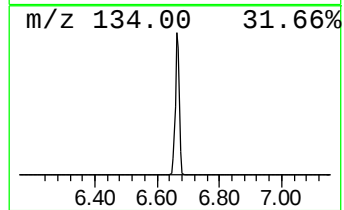
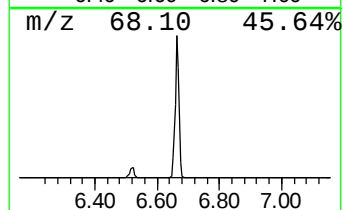
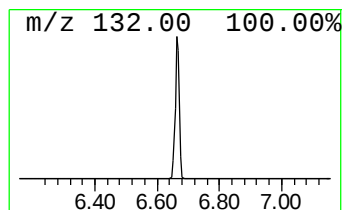
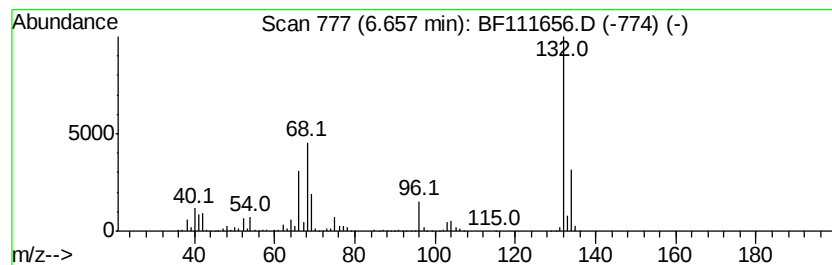
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	51.66 ng	3219980	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27	
2	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12	
3	Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10	
4	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9	
5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9	



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.12	2.6	ng	160526	1	6.90	1246650	20.0
unknown6.66	6.66	51.7	ng	3219980	1	6.90	1246650	20.0