

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF032520\
 Data File : BF119506.D
 Acq On : 25 Mar 2020 15:40
 Operator : CG/JU
 Sample : PB127810BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127810BL

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-BF031820.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	5.075	502	508	517	rBV	497637	735154	5.85%	0.896%
2	5.463	568	574	577	rBV	8397074	9725758	77.37%	11.848%
3	6.469	738	745	748	rBV	7875246	9385866	74.66%	11.434%
4	6.839	804	808	812	rVB	2558061	2018793	16.06%	2.459%
5	6.998	830	835	838	rBV	7966862	7704733	61.29%	9.386%
6	7.404	898	904	907	rBV	5676433	6463680	51.42%	7.874%
7	8.127	1021	1027	1030	rBV	2670397	2661557	21.17%	3.242%
8	9.210	1204	1211	1214	rBV	10193296	11619789	92.43%	14.155%
9	9.886	1320	1326	1329	rBV	3325895	3120344	24.82%	3.801%
10	10.674	1453	1460	1464	rBV	6844057	7689364	61.17%	9.367%
11	11.374	1572	1579	1582	rBV	3715847	3356857	26.70%	4.089%
12	12.968	1843	1850	1853	rBV	11185698	12570968	100.00%	15.314%
13	14.015	2023	2028	2031	rBV	3231914	2855501	22.72%	3.479%
14	15.480	2271	2277	2282	rBV	1723349	2180094	17.34%	2.656%

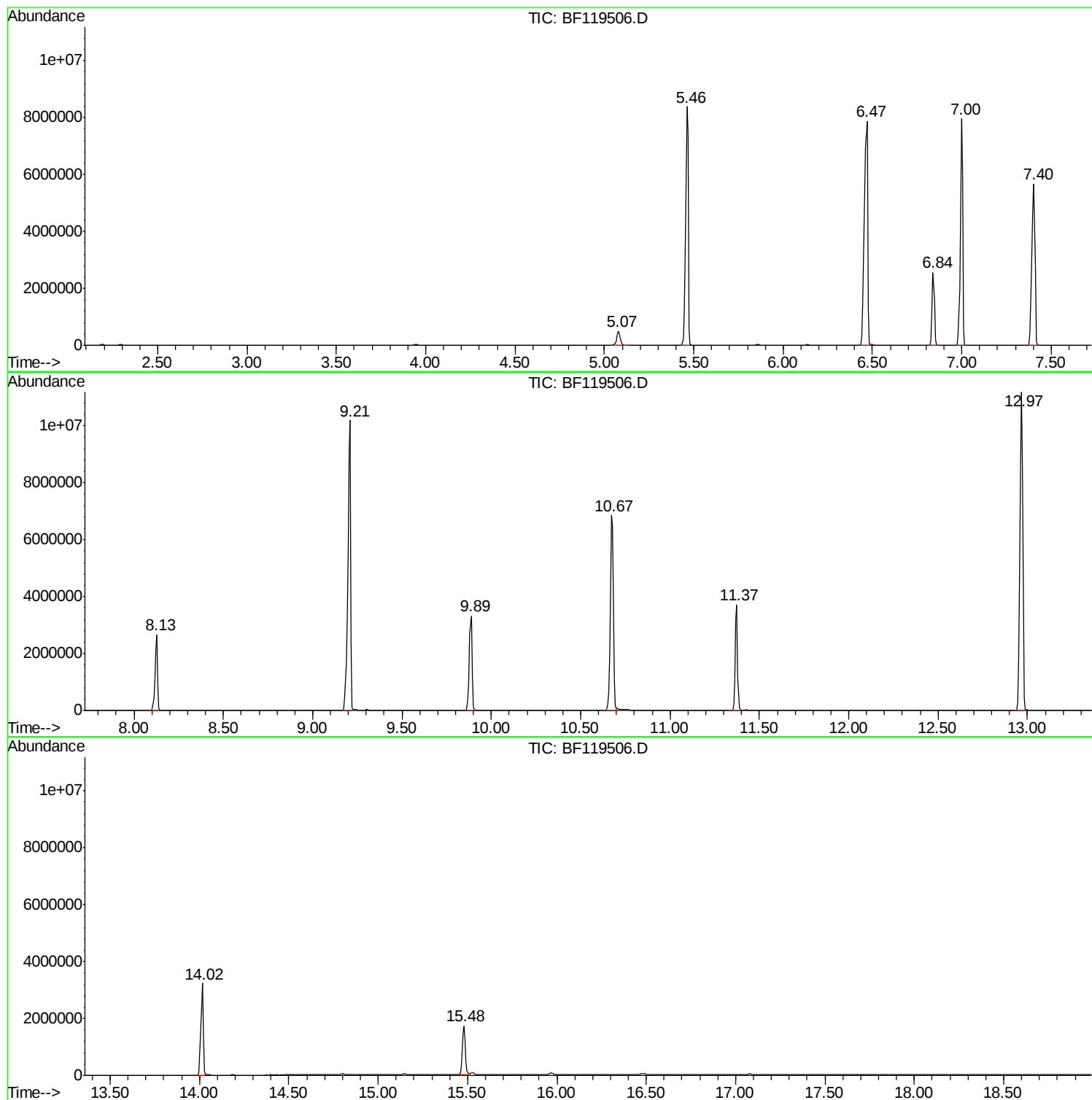
Sum of corrected areas: 82088458

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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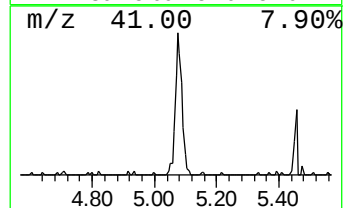
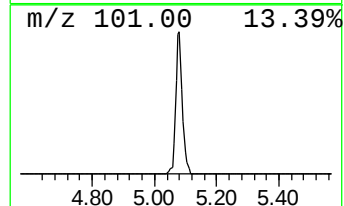
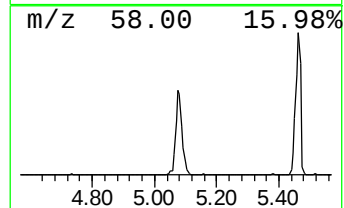
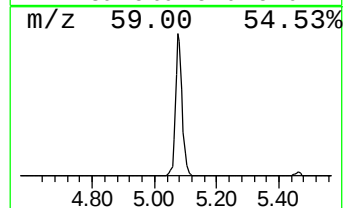
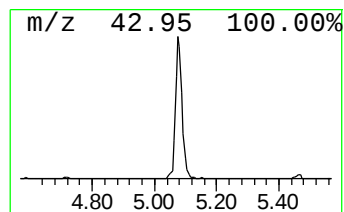
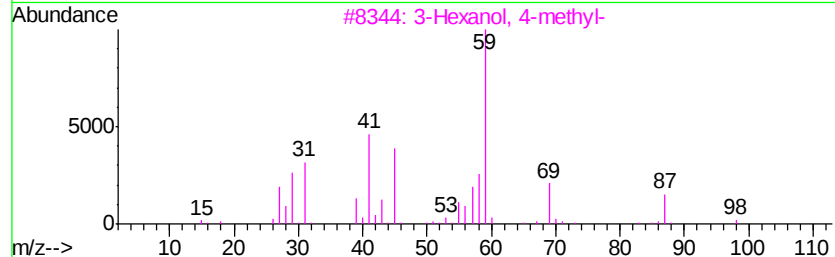
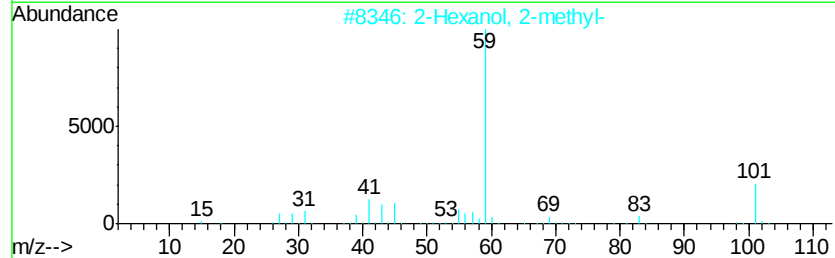
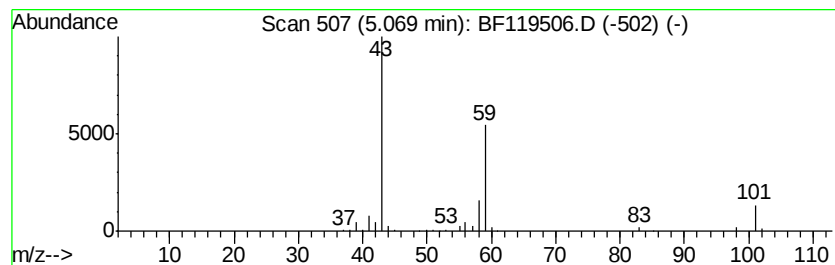
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.07	7.28 ng	735154	1,4-Dichlorobenzene-d4	6.84

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	33
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			3-Octanol	130	C8H18O	000589-98-0	33
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	7.3	ng	735154	1	6.84	2018790	20.0