

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF073118\
 Data File : BF107870.D
 Acq On : 1 Aug 2018 00:34
 Operator : JU/SJ
 Sample : J3825-13 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-073018-B

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF073018.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	5.184	481	484	495	rBV	209794	247654	13.20%	1.453%
2	5.595	551	554	579	rBV	568950	1198800	63.91%	7.035%
3	6.601	722	725	744	rBV	546360	1296630	69.12%	7.609%
4	6.736	744	748	783	rVV	1096996	1875779	100.00%	11.008%
5	6.960	783	786	809	rVV	590497	1047507	55.84%	6.147%
6	7.113	809	812	834	rVB	990869	1249857	66.63%	7.335%
7	7.530	879	883	905	rBV	285654	819860	43.71%	4.811%
8	8.242	1001	1004	1025	rBV	863598	1250637	66.67%	7.340%
9	9.313	1183	1186	1215	rBV	1142816	1651924	88.07%	9.695%
10	9.713	1249	1254	1266	rVB	51534	79681	4.25%	0.468%
11	10.001	1299	1303	1315	rBV	980907	1120288	59.72%	6.575%
12	10.801	1435	1439	1453	rBV	422926	753942	40.19%	4.425%
13	11.501	1554	1558	1571	rBV	672603	1033917	55.12%	6.068%
14	12.001	1640	1643	1646	rBV5	13386	21419	1.14%	0.126%
15	12.718	1763	1765	1769	rBV	28796	41667	2.22%	0.245%
16	12.954	1802	1805	1810	rVB2	50945	67997	3.63%	0.399%
17	13.077	1823	1826	1837	rBV	1120914	1430813	76.28%	8.397%
18	14.154	2005	2009	2019	rBV	810861	1088478	58.03%	6.388%
19	15.695	2268	2271	2284	rVB	417948	762911	40.67%	4.477%

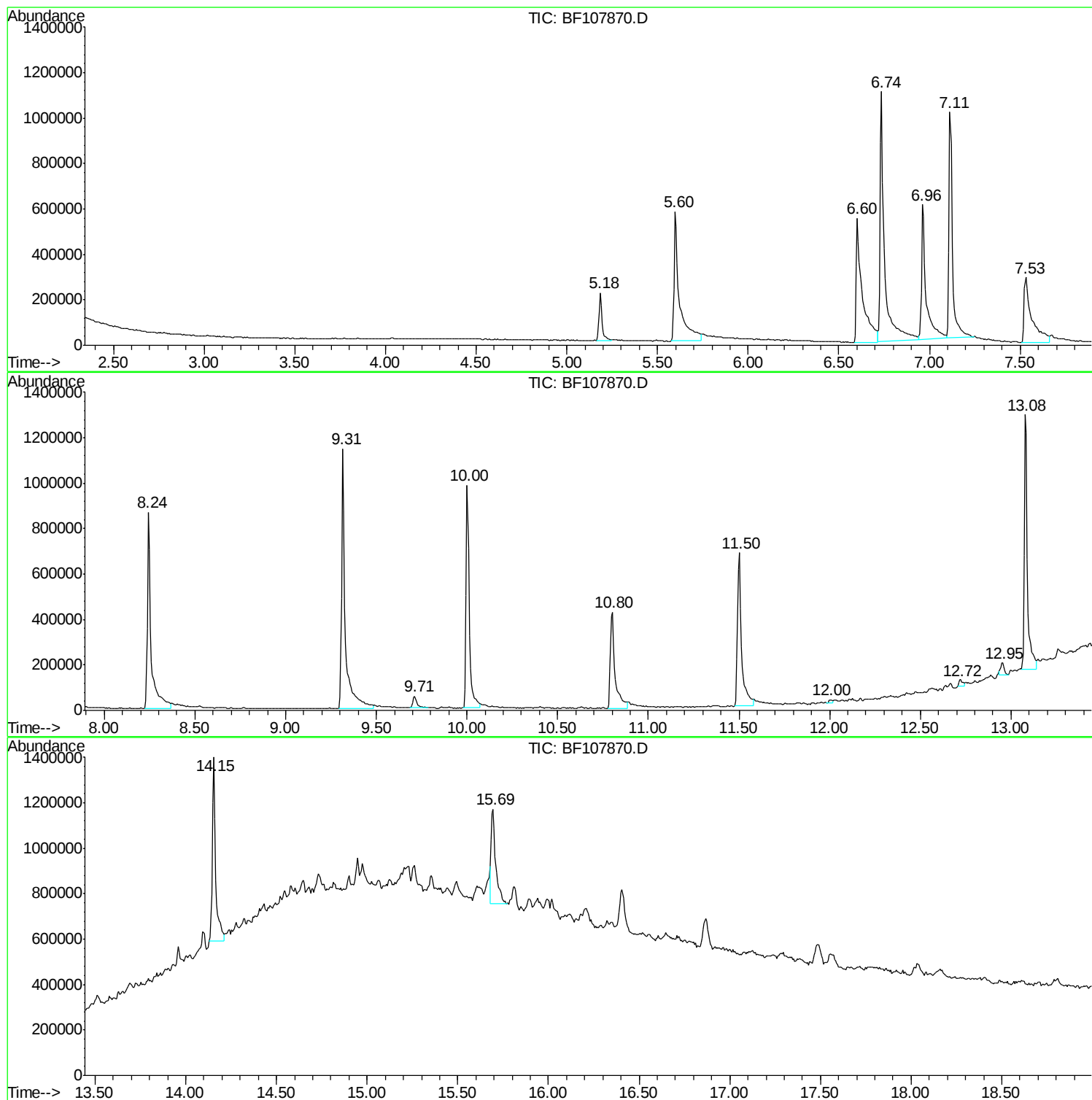
Sum of corrected areas: 17039761

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.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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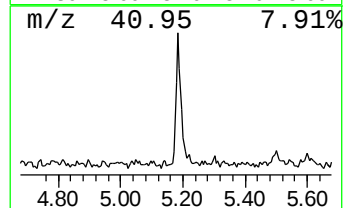
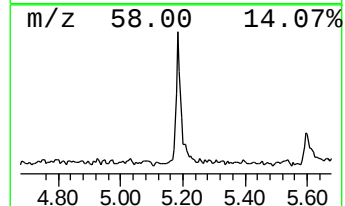
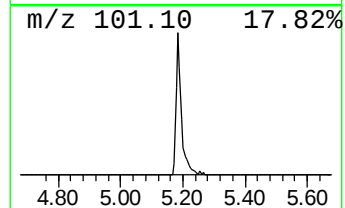
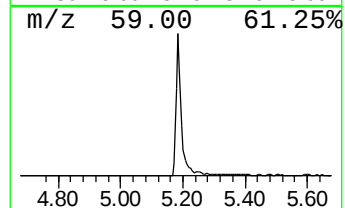
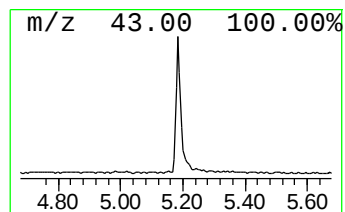
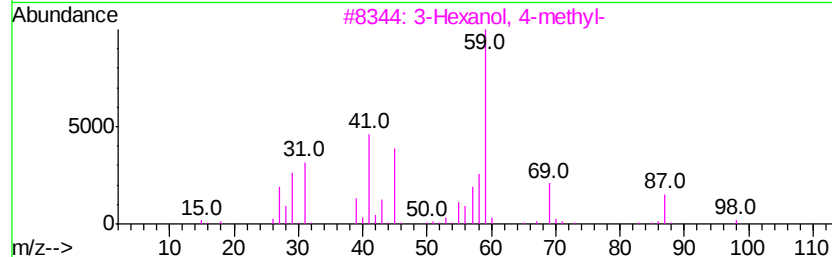
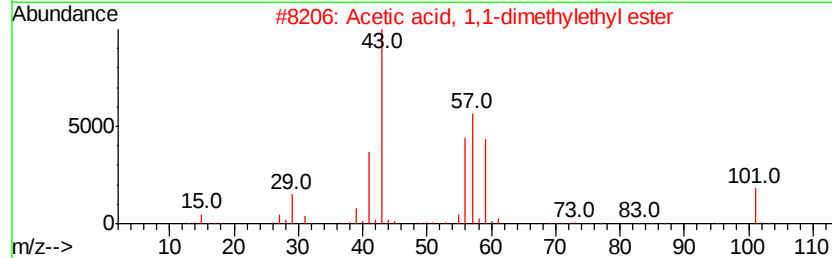
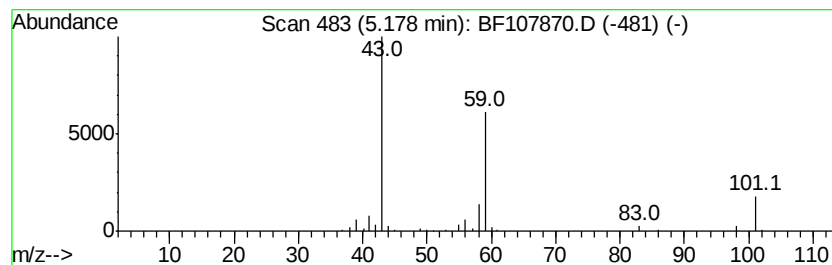
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.18	4.73 ng	247654	1,4-Dichlorobenzene-d4	6.96

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
4			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25



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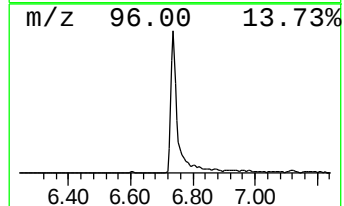
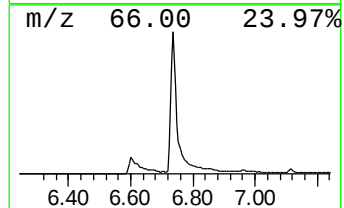
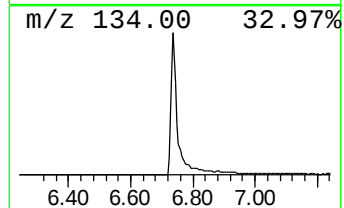
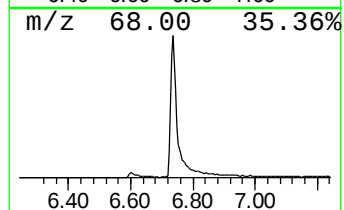
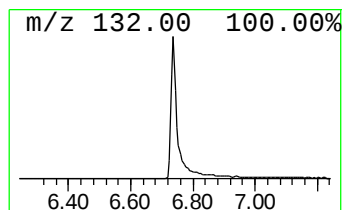
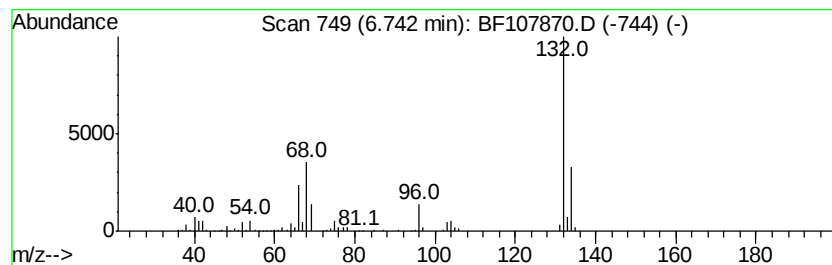
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	35.81 ng	1875780	1,4-Dichlorobenzene-d4	6.96

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	35
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.18	4.7	ng	247654	1	6.96	1047510	20.0
unknown6.74	6.74	35.8	ng	1875780	1	6.96	1047510	20.0