

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF122819\
 Data File : BF118364.D
 Acq On : 28 Dec 2019 14:37
 Operator : JU
 Sample : PB125749BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB125749BL

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-BF122419.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	5.081	504	509	525	rBV	478752	685374	11.20%	1.702%
2	5.487	573	578	600	rBV	3707235	3771616	61.62%	9.366%
3	6.493	743	749	752	rBV	3656620	3781867	61.79%	9.391%
4	6.628	767	772	775	rBV	4547149	4229683	69.10%	10.504%
5	6.851	806	810	818	rBV	1018598	841697	13.75%	2.090%
6	7.010	832	837	840	rBV	3757116	3345925	54.66%	8.309%
7	7.422	901	907	922	rBV	2113136	2510835	41.02%	6.235%
8	8.140	1024	1029	1044	rBV	1068547	1093176	17.86%	2.715%
9	9.216	1206	1212	1227	rBV	4977503	5366222	87.67%	13.326%
10	9.898	1323	1328	1340	rVB	1331648	1337341	21.85%	3.321%
11	10.698	1459	1464	1482	rBV	3715094	3586314	58.59%	8.906%
12	11.398	1579	1583	1600	rVB	1354618	1426353	23.30%	3.542%
13	12.992	1848	1854	1857	rBV	5944624	6120967	100.00%	15.200%
14	14.051	2030	2034	2042	rBV	1133633	1108273	18.11%	2.752%
15	15.151	2217	2221	2226	rBV	56253	64493	1.05%	0.160%
16	15.545	2282	2288	2302	rBV	591944	927958	15.16%	2.304%
17	15.980	2358	2362	2367	rBV2	49250	71165	1.16%	0.177%

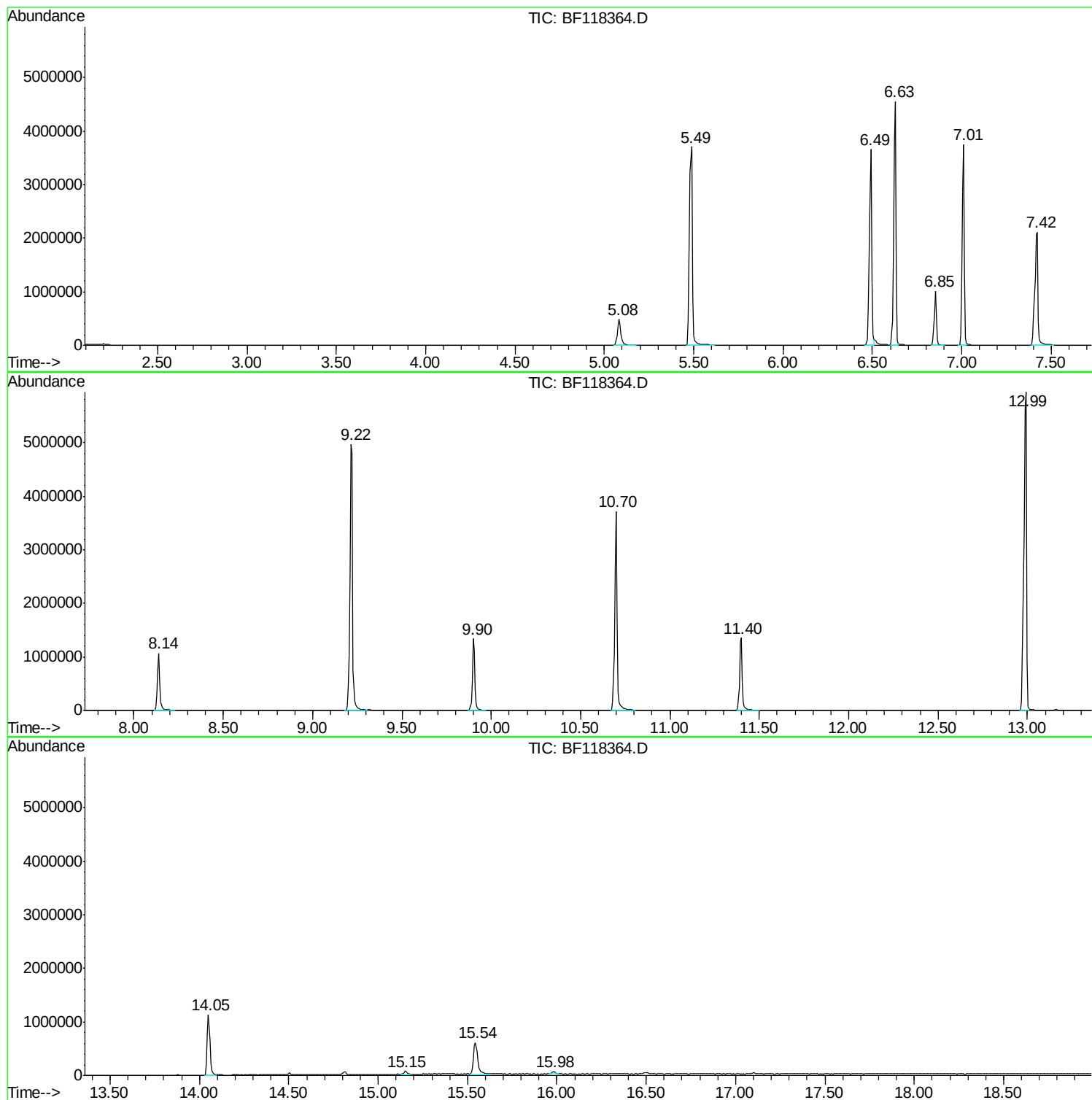
Sum of corrected areas: 40269259

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF122419.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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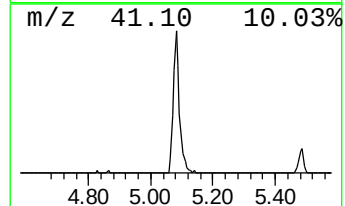
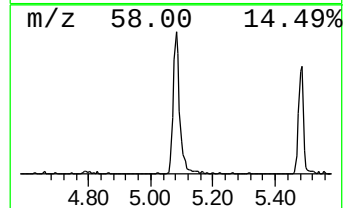
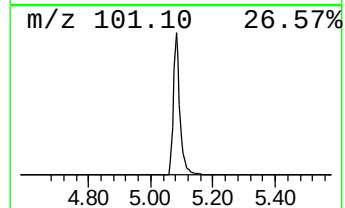
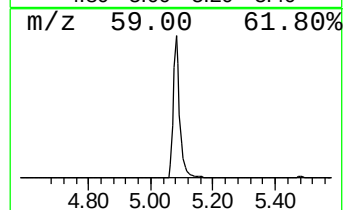
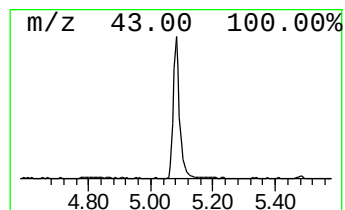
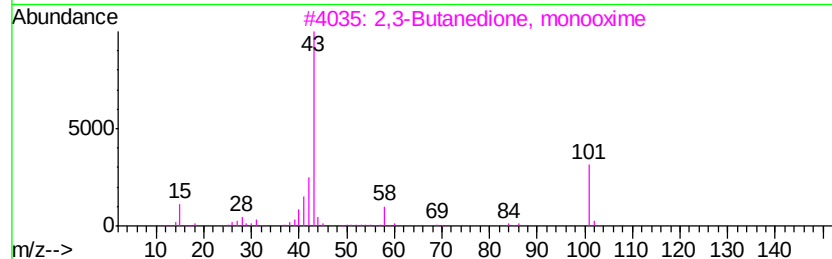
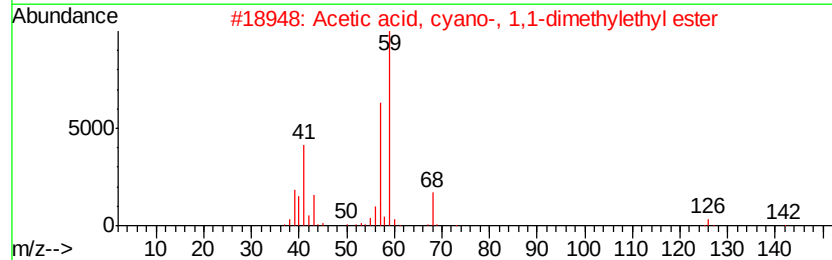
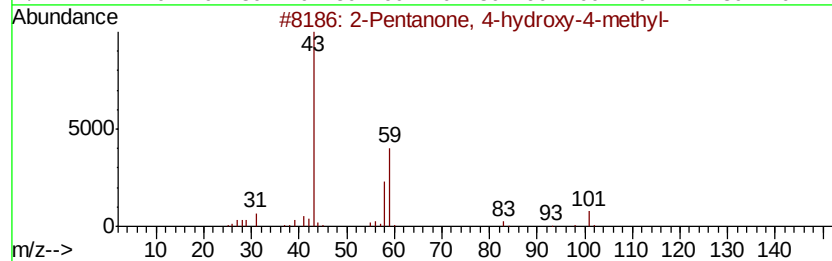
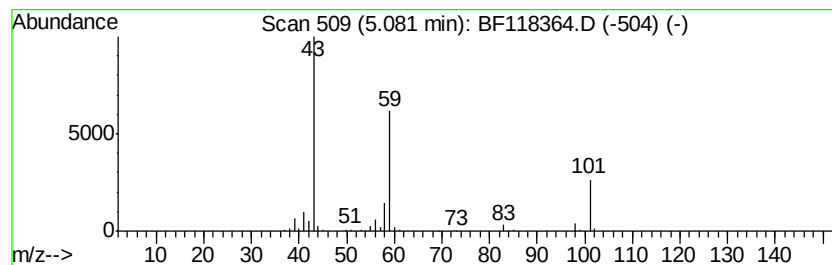
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.08	16.29 ng	685374	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4		5-Hexen-2-one	98	C6H10O	000109-49-9	9
5		Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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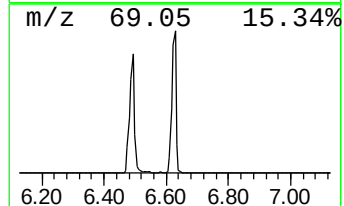
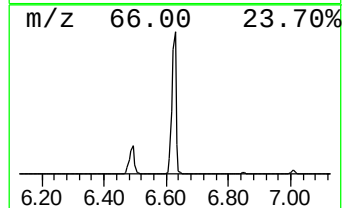
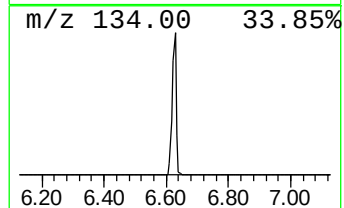
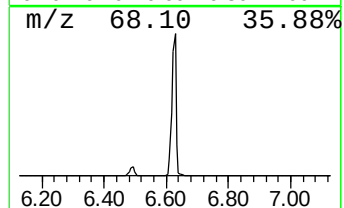
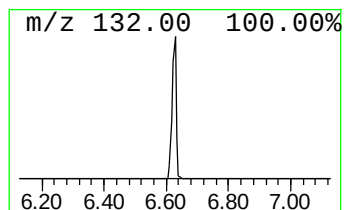
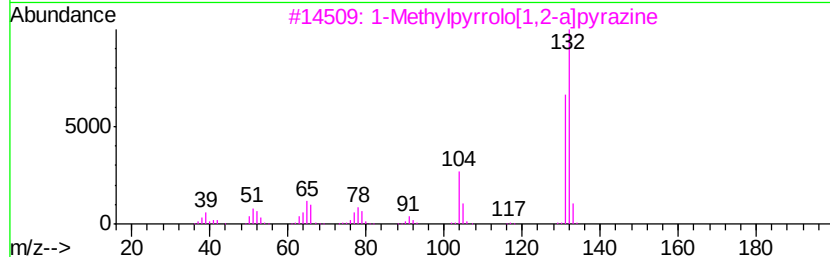
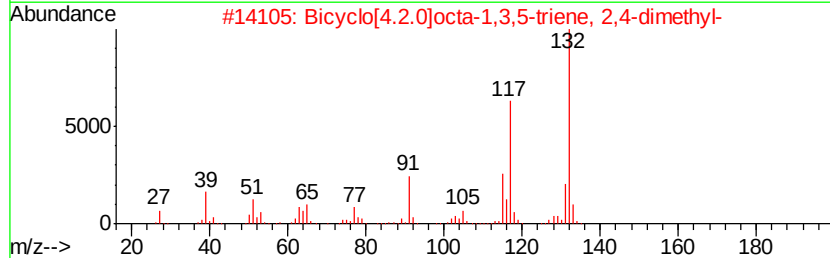
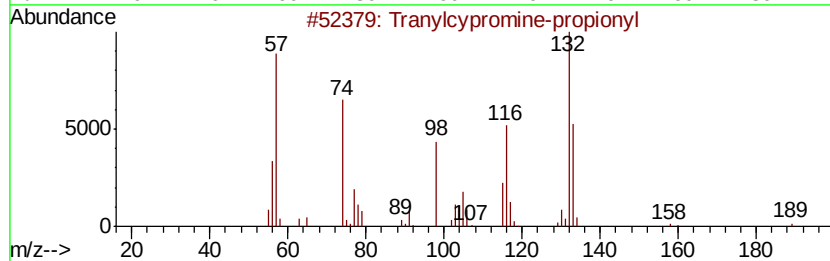
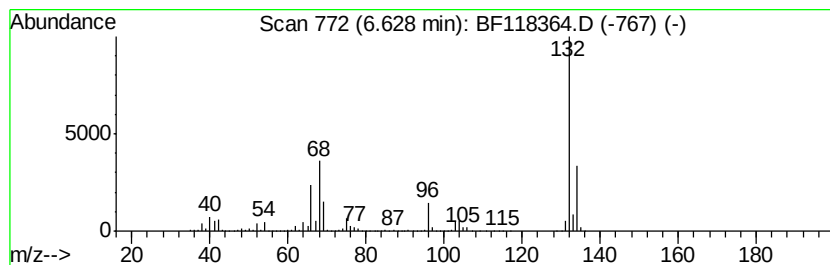
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Peak Number 2 unknown6.63 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	100.50 ng	4229680	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	16
2		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
3		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
4		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.08	16.3	ng	685374	1	6.85	841697	20.0
unknown6.63	6.63	100.5	ng	4229680	1	6.85	841697	20.0