

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072018\
 Data File : BG035789.D
 Acq On : 20 Jul 2018 18:09
 Operator : JU/SJ
 Sample : PB111269BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB111269BL

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_G\METHODS\8270-BG071218.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	5.138	310	315	326	rBV	124079	200659	4.63%	1.027%
2	5.614	388	396	412	rBV	876858	1508025	34.82%	7.716%
3	7.200	658	666	679	rBV	885569	1609532	37.17%	8.235%
4	7.564	720	728	740	rBV	900888	1547606	35.74%	7.919%
5	8.029	801	807	816	rVB	175430	304637	7.03%	1.559%
6	8.352	853	862	871	rBV	682192	1212991	28.01%	6.207%
7	9.192	996	1005	1021	rBV	502828	969984	22.40%	4.963%
8	10.831	1278	1284	1295	rBV	183161	375039	8.66%	1.919%
9	13.274	1692	1700	1714	rBV	1504130	2371137	54.75%	12.132%
10	14.655	1928	1935	1949	rBV2	321112	566546	13.08%	2.899%
11	16.141	2181	2188	2203	rBV	1139537	1880466	43.42%	9.622%
12	17.398	2394	2402	2419	rBV	463425	805982	18.61%	4.124%
13	20.001	2839	2845	2860	rBV	3234045	4330631	100.00%	22.159%
14	21.657	3120	3127	3137	rBV	547067	985839	22.76%	5.044%
15	24.859	3661	3672	3684	rBV	292512	874779	20.20%	4.476%

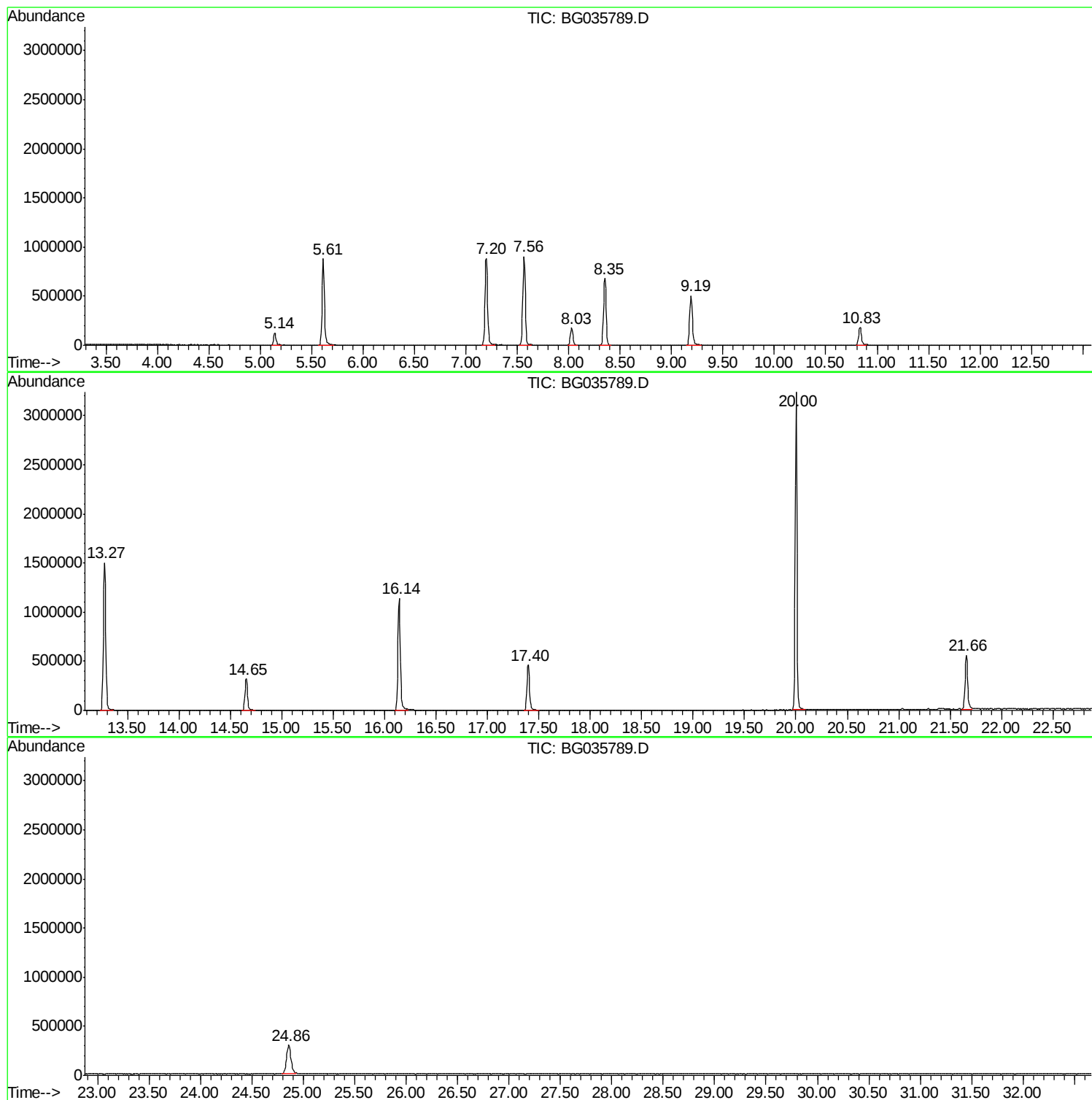
Sum of corrected areas: 19543853

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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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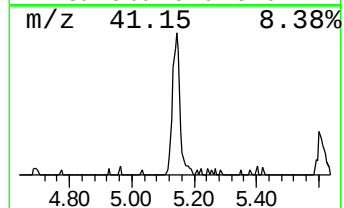
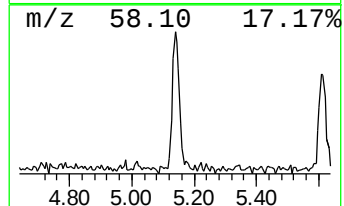
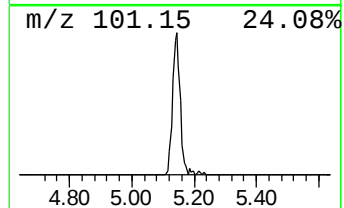
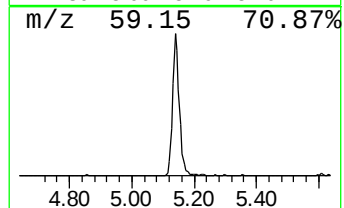
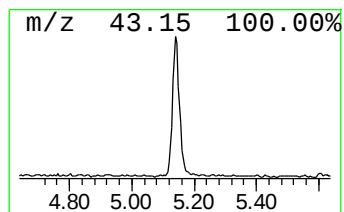
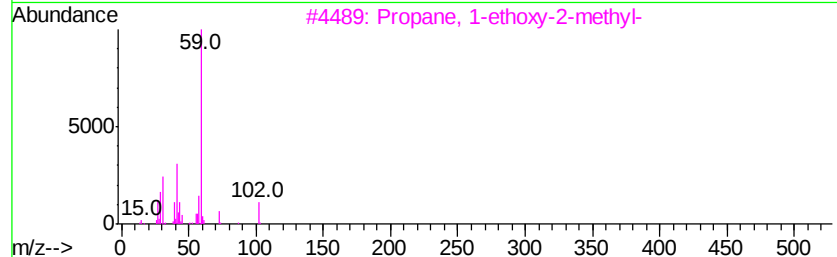
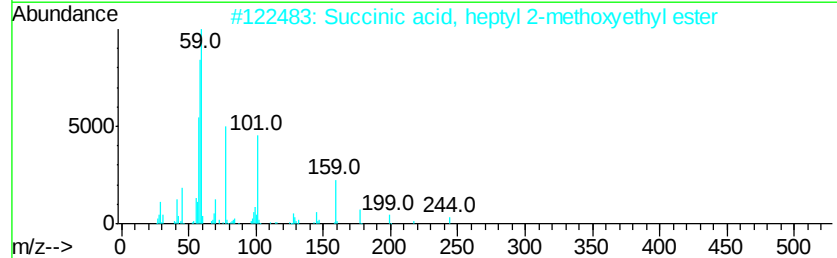
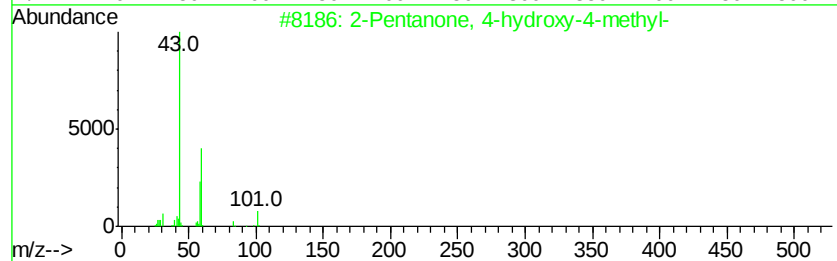
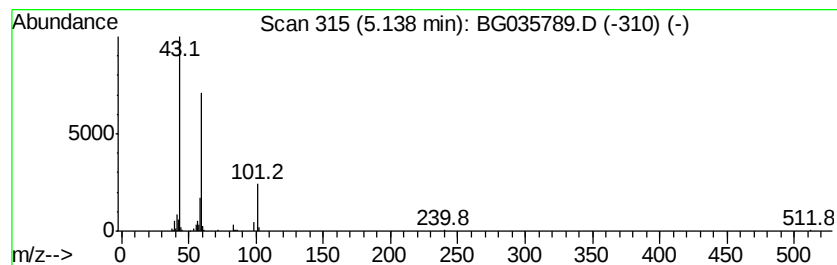
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	13.17 ng	200659	1,4-Dichlorobenzene-d4	8.03

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Succinic acid, heptyl 2-methoxyethyl ester	274	C14H26O5	1000325-80-7	33
3			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
4			Hydrazine, 1,1-bis(1-methylethyl)-	116	C6H16N2	000921-14-2	25
5			Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	17



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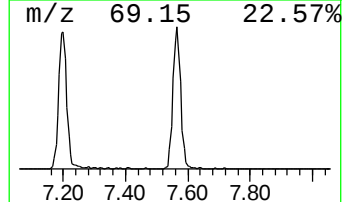
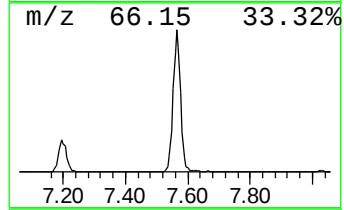
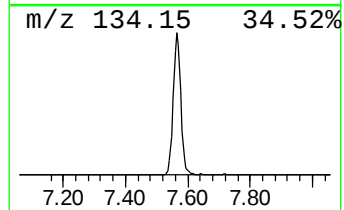
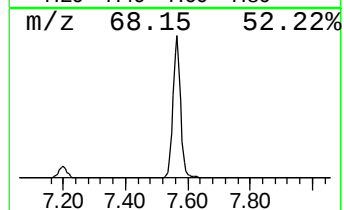
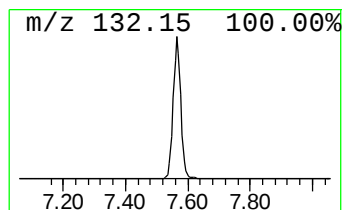
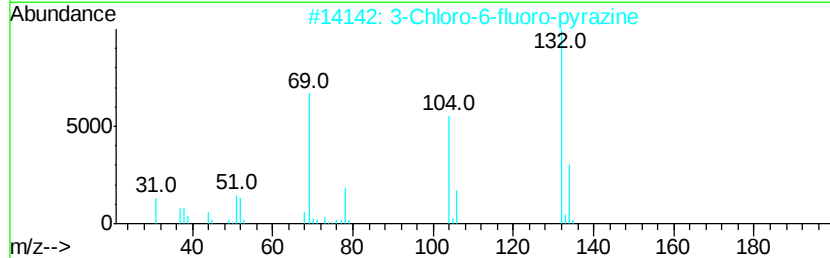
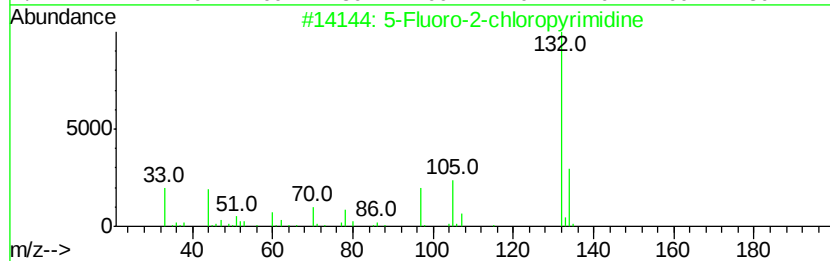
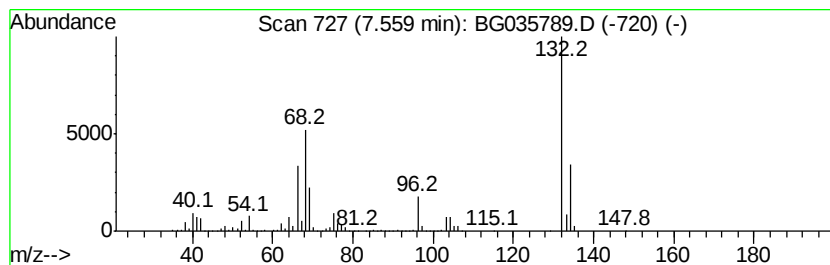
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	101.60 ng	1547610	1,4-Dichlorobenzene-d4	8.03

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	27
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
3		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	10
4		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10
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...	5.14	13.2	ng	200659	1	8.03	304637	20.0
unknown7.56	7.56	101.6	ng	1547610	1	8.03	304637	20.0