

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG042419\
 Data File : BG040601.D
 Acq On : 24 Apr 2019 19:49
 Operator : JU/SJ
 Sample : PB119187BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB119187BL

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_G\METHODS\8270-BG042319.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	3.159	7	12	21	rVB3	41693	85775	1.86%	0.418%
2	5.685	434	442	451	rBV	1005460	1548410	33.55%	7.548%
3	7.295	705	716	724	rBV	1024515	1774095	38.44%	8.648%
4	7.677	773	781	792	rBV	946120	1654991	35.86%	8.068%
5	8.159	855	863	869	rBV	188465	320180	6.94%	1.561%
6	8.482	908	918	927	rBV	681461	1166730	25.28%	5.688%
7	9.334	1055	1063	1071	rBV	595060	1045663	22.66%	5.097%
8	10.979	1336	1343	1351	rVB	249904	434961	9.42%	2.120%
9	13.411	1750	1757	1765	rVB	1542704	2363105	51.20%	11.520%
10	14.786	1983	1991	1999	rBV2	406109	646301	14.00%	3.151%
11	16.267	2235	2243	2252	rBV	1082827	1664972	36.07%	8.116%
12	17.530	2452	2458	2468	rVB	593438	889096	19.26%	4.334%
13	20.127	2894	2900	2906	rVB	3473534	4615591	100.00%	22.500%
14	21.819	3181	3188	3198	rVB	692665	1163551	25.21%	5.672%
15	25.192	3751	3762	3775	rVB3	374147	1140087	24.70%	5.558%

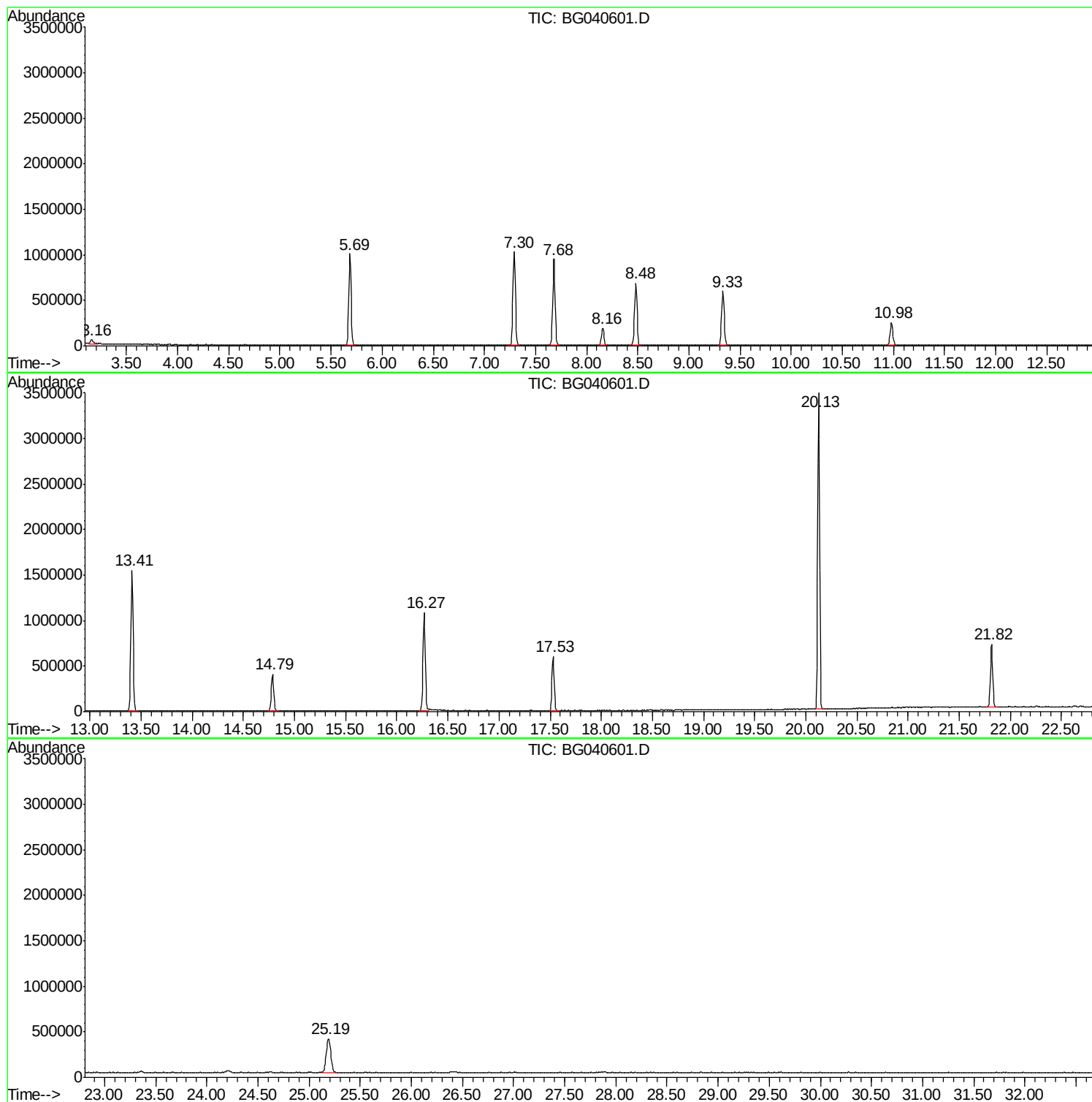
Sum of corrected areas: 20513508

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



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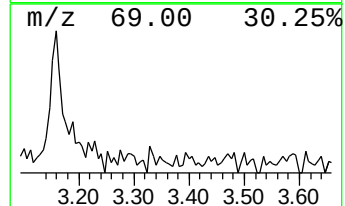
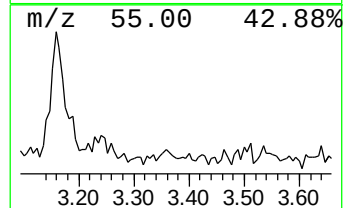
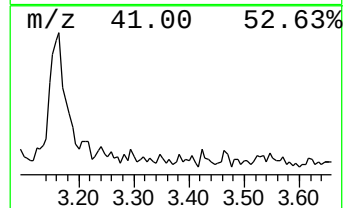
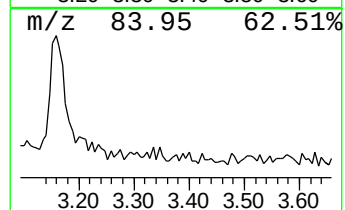
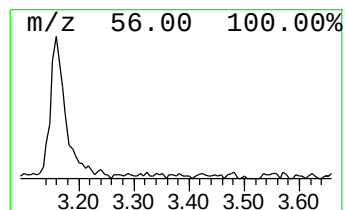
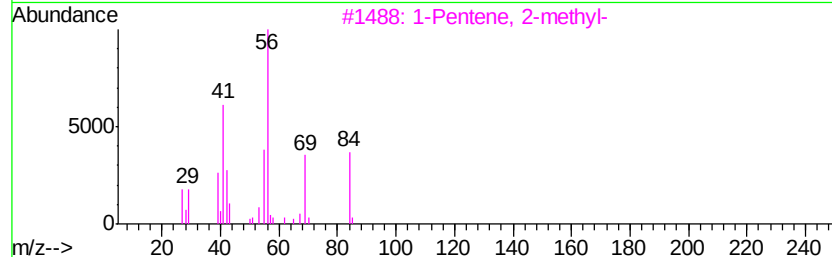
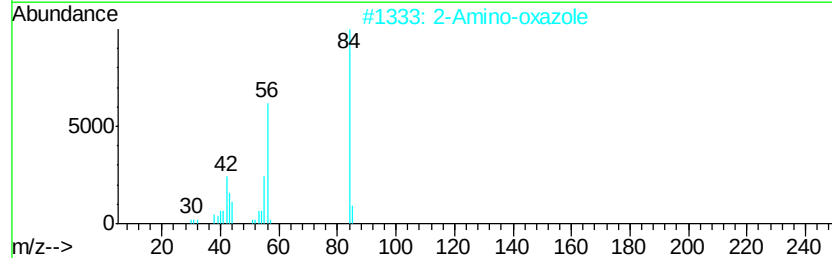
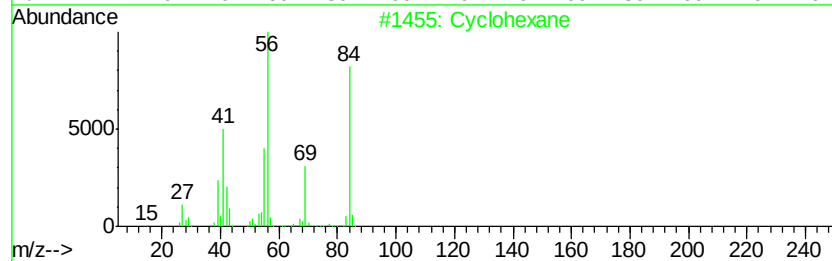
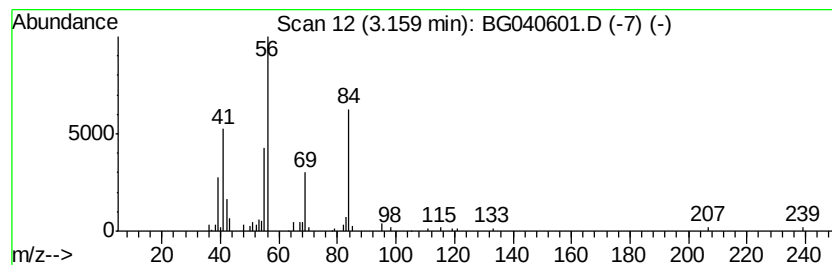
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Peak Number 1 1-Pentene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.16	5.36 ng	85775	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclohexane	84	C6H12	000110-82-7	87
2		2-Amino-oxazole	84	C3H4N2O	004570-45-0	64
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	56
4		1-Heptene, 2-methyl-	112	C8H16	015870-10-7	50
5		Iron pentacarbonyl	196	C5FeO5	013463-40-6	50



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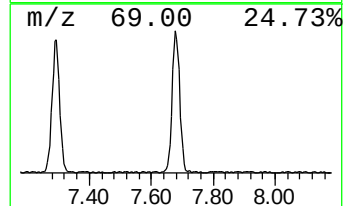
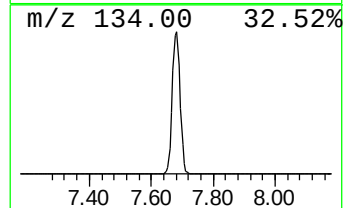
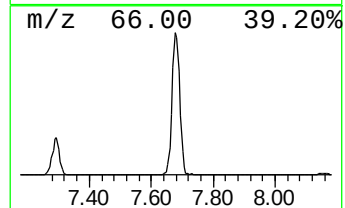
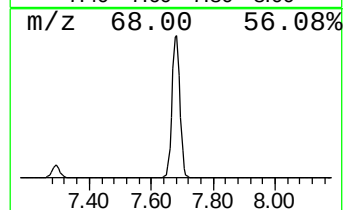
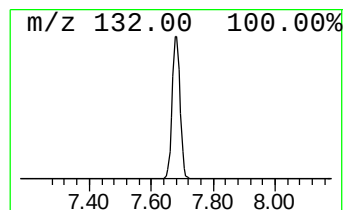
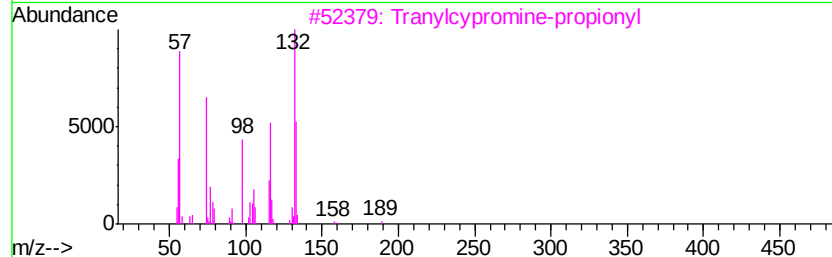
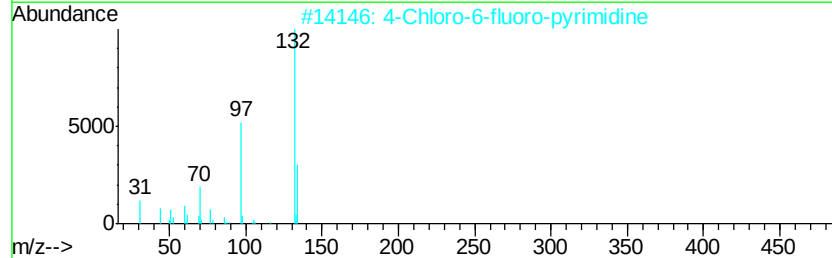
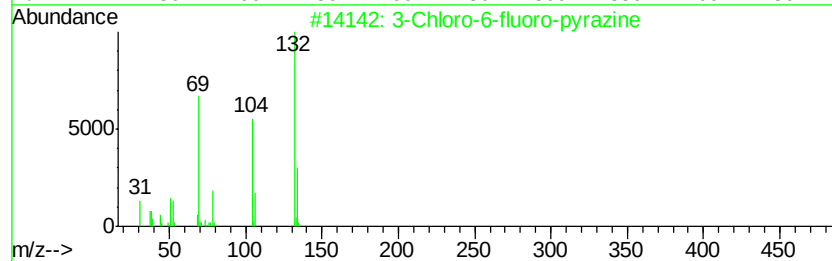
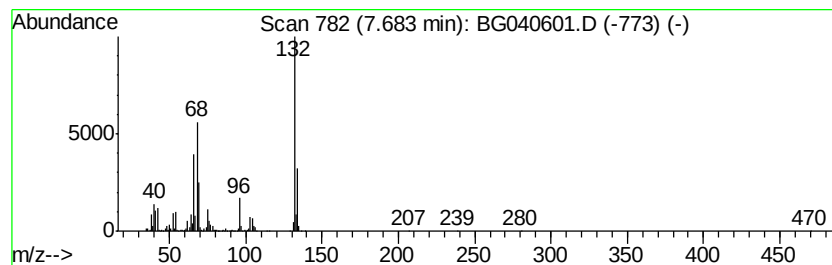
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Peak Number 2 unknown7.68 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.68	103.38 ng	1654990	1,4-Dichlorobenzene-d4	8.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
2		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	14
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	14
4		1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	12
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Pentene, 2-methyl-	3.16	5.4	ng	85775	1	8.16	320180	20.0
unknown7.68	7.68	103.4	ng	1654990	1	8.16	320180	20.0