

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG072718\
 Data File : BG035942.D
 Acq On : 27 Jul 2018 18:02
 Operator : JU/SJ
 Sample : J4192-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

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-BG071218.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

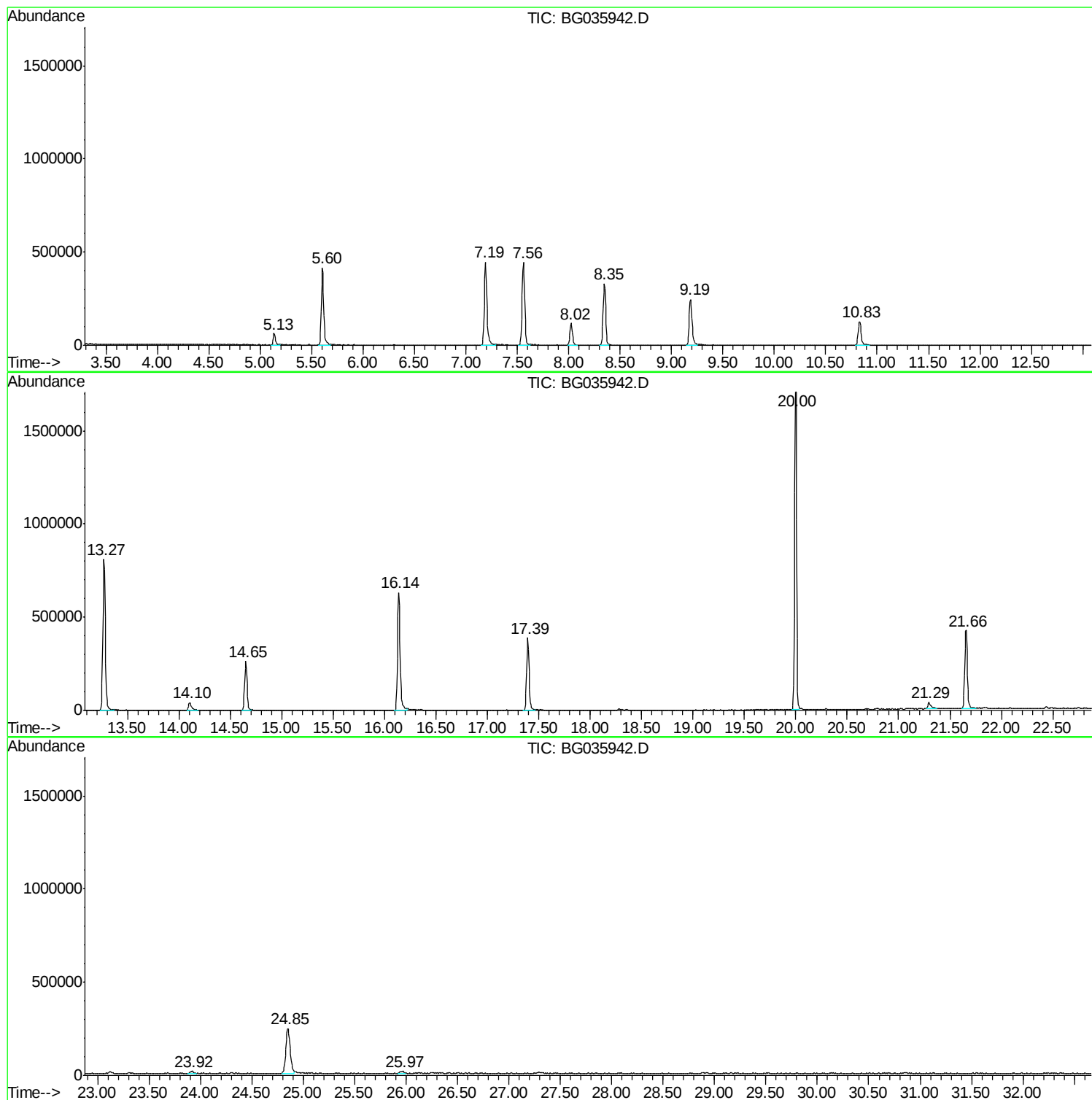
peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.133	309	314	324	rBV	61372	97528	4.09%	0.846%
2	5.603	388	394	408	rBV	415111	716797	30.08%	6.217%
3	7.189	657	664	681	rBV	446978	845500	35.48%	7.333%
4	7.559	719	727	738	rBV	443221	779137	32.70%	6.758%
5	8.023	800	806	816	rBV	118668	196816	8.26%	1.707%
6	8.346	853	861	869	rBV	332338	590921	24.80%	5.125%
7	9.186	996	1004	1015	rBV	245351	496169	20.82%	4.304%
8	10.825	1276	1283	1299	rBV	126468	265602	11.15%	2.304%
9	13.269	1692	1699	1716	rBV	809046	1315896	55.22%	11.414%
10	14.103	1836	1841	1854	rBV2	41631	89119	3.74%	0.773%
11	14.649	1927	1934	1943	rBV2	264066	434674	18.24%	3.770%
12	16.136	2180	2187	2201	rBV	631963	1078847	45.28%	9.357%
13	17.393	2393	2401	2413	rBV	389556	640241	26.87%	5.553%
14	19.995	2838	2844	2851	rBV	1704501	2382845	100.00%	20.668%
15	21.293	3060	3065	3077	rBV6	33931	69703	2.93%	0.605%
16	21.658	3119	3127	3143	rBV2	420490	758615	31.84%	6.580%
17	23.919	3505	3512	3518	rVB5	13239	26369	1.11%	0.229%
18	24.853	3660	3671	3681	rBV2	240074	710725	29.83%	6.165%
19	25.969	3852	3861	3865	rBV8	14375	33785	1.42%	0.293%

Sum of corrected areas: 11529289

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG071218.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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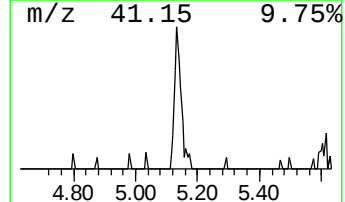
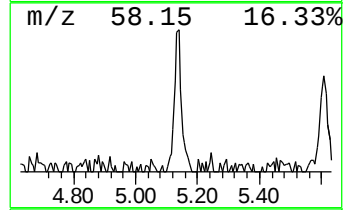
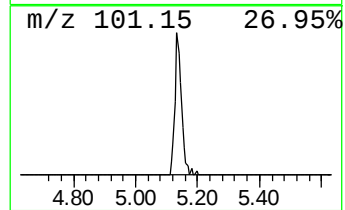
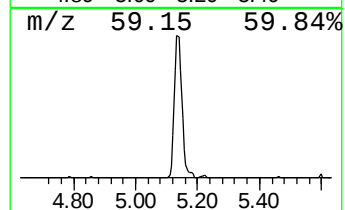
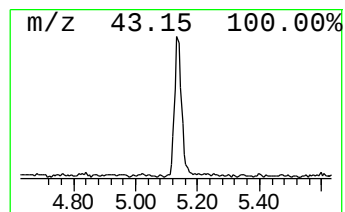
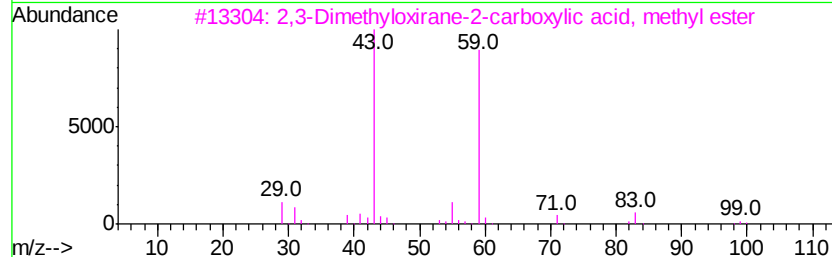
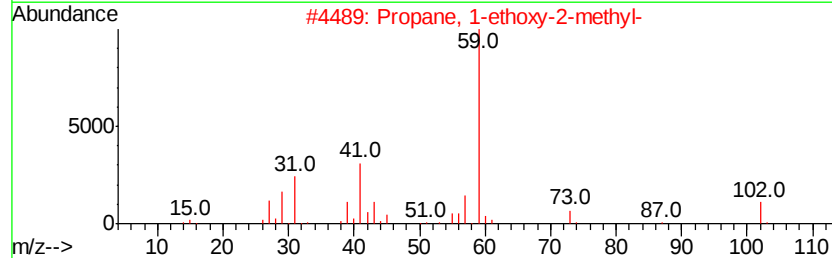
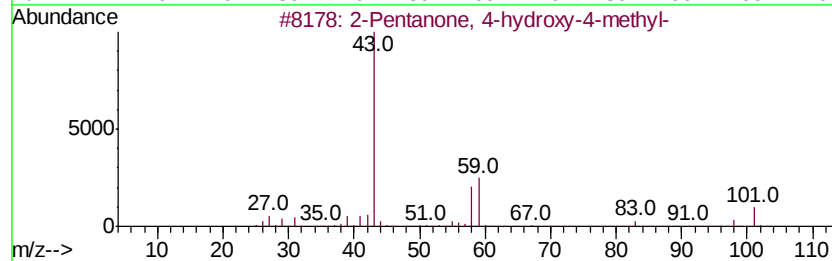
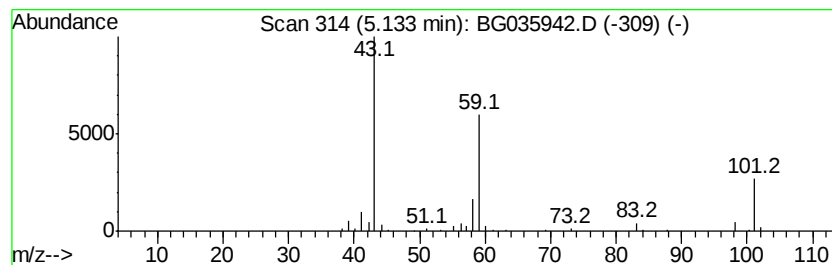
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	9.91 ng	97528	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2			Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	38
3			2,3-Dimethyloxirane-2-carboxylic...	130	C6H10O3	1000306-64-4	28
4			2-Pentanol, 2,4-dimethyl-	116	C7H16O	000625-06-9	25
5			3-Hexanol	102	C6H14O	000623-37-0	25



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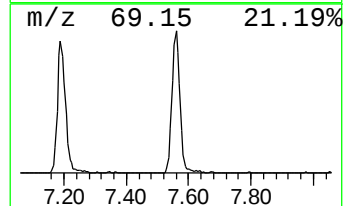
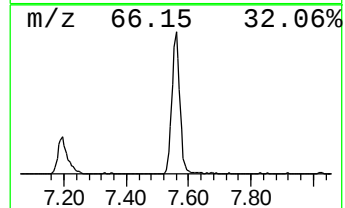
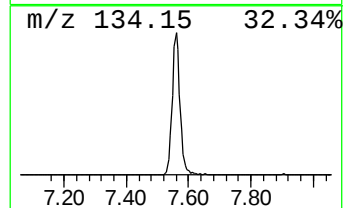
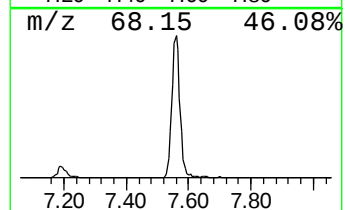
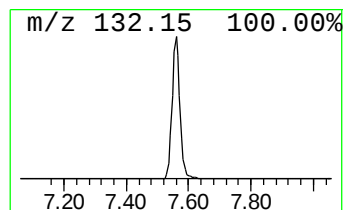
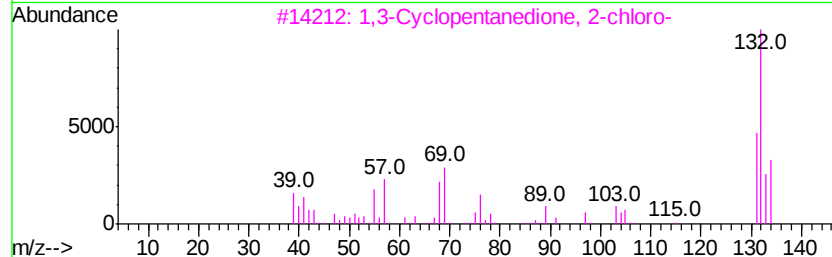
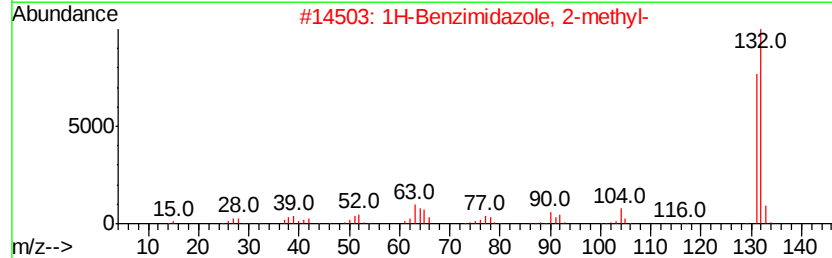
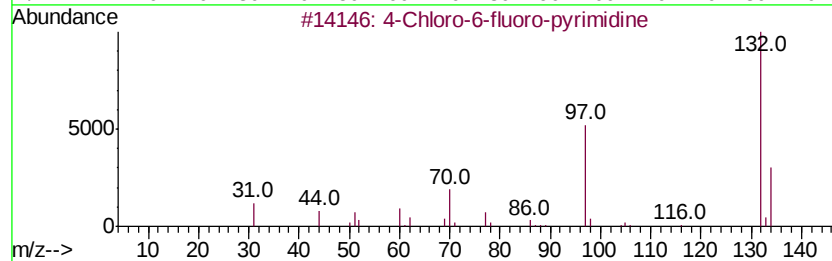
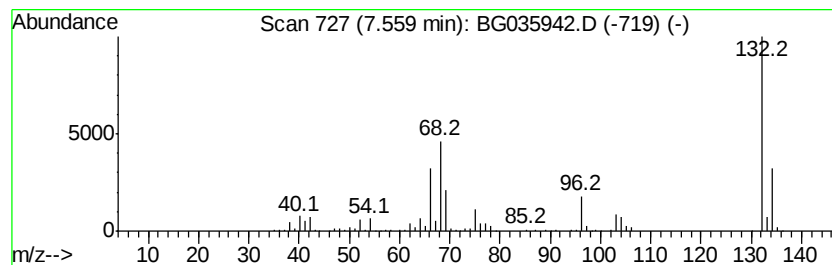
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Peak Number 2 unknown7.56 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.56	79.17 ng	779137	1,4-Dichlorobenzene-d4	8.02

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
3			1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17
4			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5			Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	9.9	ng	97528	1	8.02	196816	20.0
unknown7.56	7.56	79.2	ng	779137	1	8.02	196816	20.0