

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010719\
 Data File : BG038983.D
 Acq On : 8 Jan 2019 1:43
 Operator : JU/SJ
 Sample : PB116227BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB116227BL

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-BG121218.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	5.132	342	348	357	rBV	29251	46631	2.07%	0.421%
2	5.602	421	428	442	rBV	499826	842483	37.38%	7.603%
3	7.212	693	702	714	rBV	521505	928589	41.21%	8.380%
4	7.576	755	764	777	rBV	504942	895956	39.76%	8.086%
5	8.046	836	844	850	rBV	104312	182354	8.09%	1.646%
6	8.369	890	899	907	rBV	388165	698096	30.98%	6.300%
7	9.221	1037	1044	1054	rBV	301132	561014	24.89%	5.063%
8	10.860	1312	1323	1334	rBV	130734	246633	10.94%	2.226%
9	13.299	1730	1738	1748	rBV	870258	1448511	64.28%	13.073%
10	14.685	1967	1974	1985	rBV2	218988	370911	16.46%	3.347%
11	16.178	2221	2228	2242	rBV	722747	1174885	52.13%	10.603%
12	17.429	2435	2441	2453	rBV	291245	471831	20.94%	4.258%
13	20.032	2878	2884	2894	rBV	1685113	2253579	100.00%	20.338%
14	21.706	3163	3169	3177	rBV	287402	490298	21.76%	4.425%
15	24.997	3719	3729	3742	rVB2	153104	468606	20.79%	4.229%

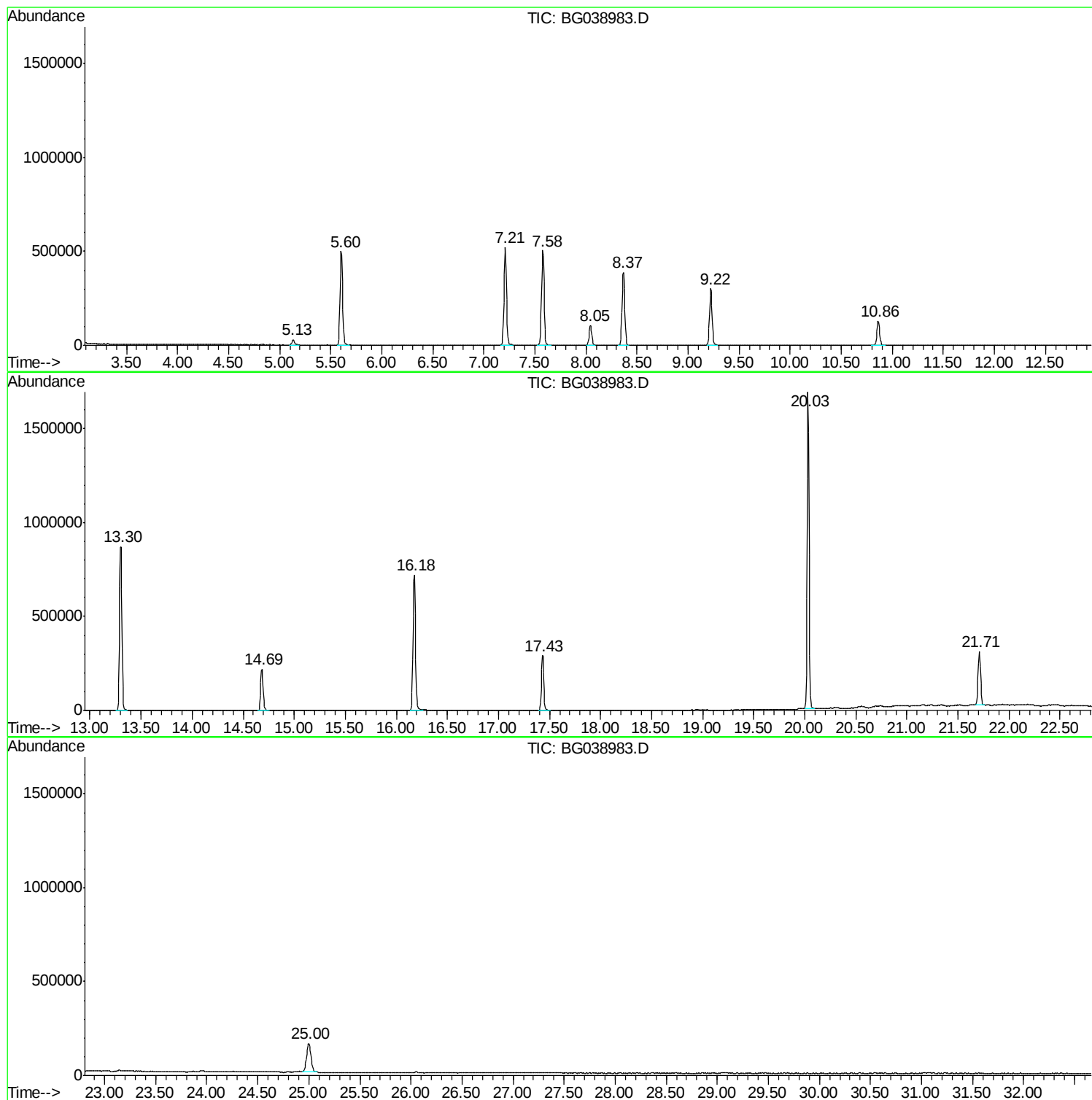
Sum of corrected areas: 11080377

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



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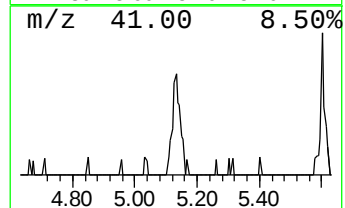
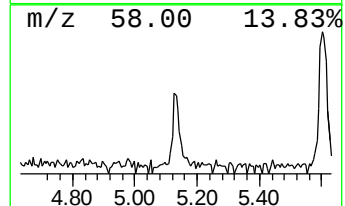
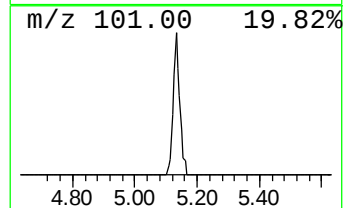
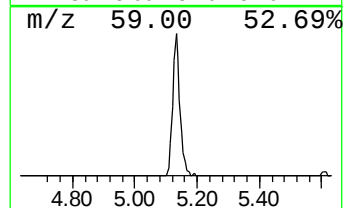
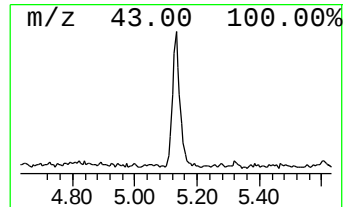
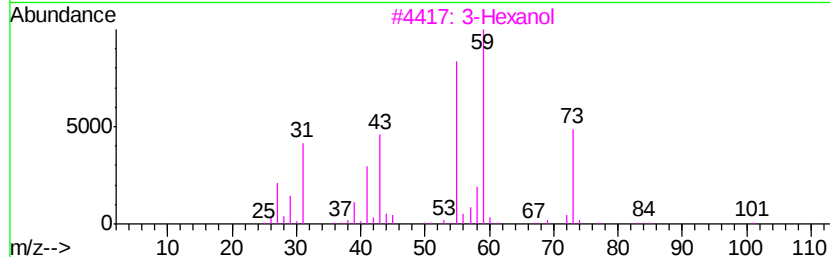
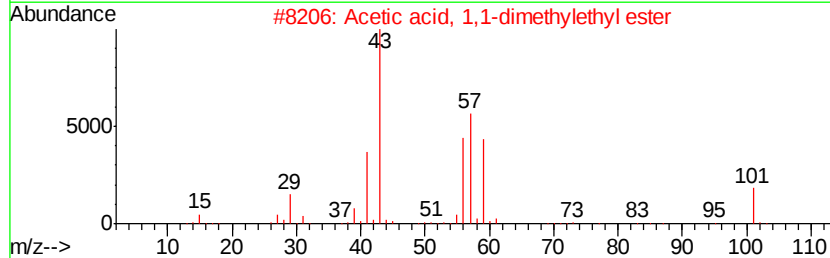
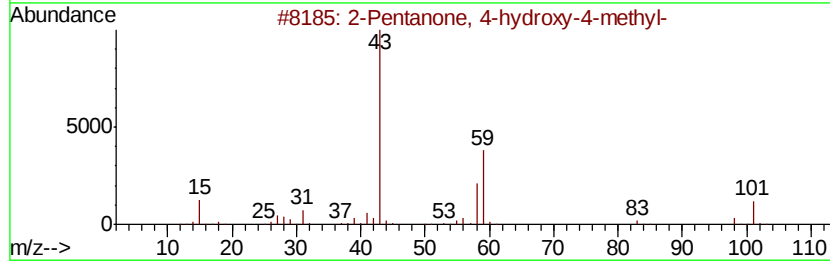
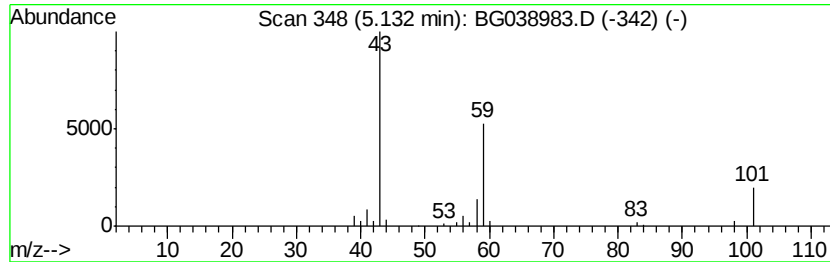
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.13	5.11 ng	46631	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3			3-Hexanol	102	C6H14O	000623-37-0	25
4			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
5			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17



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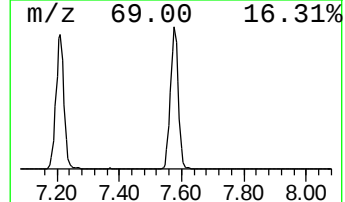
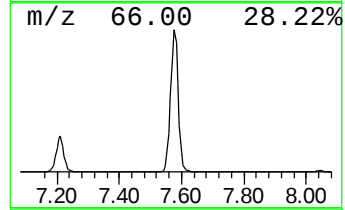
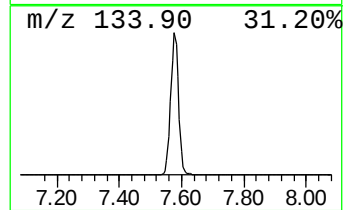
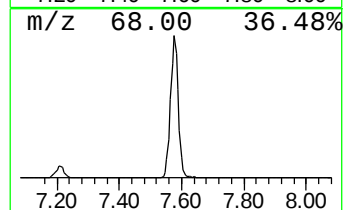
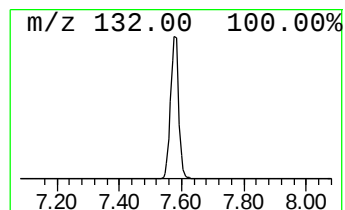
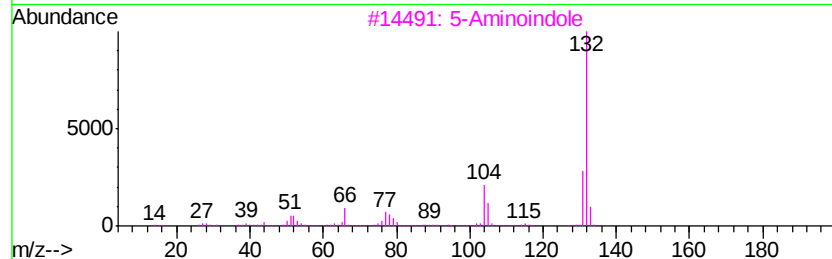
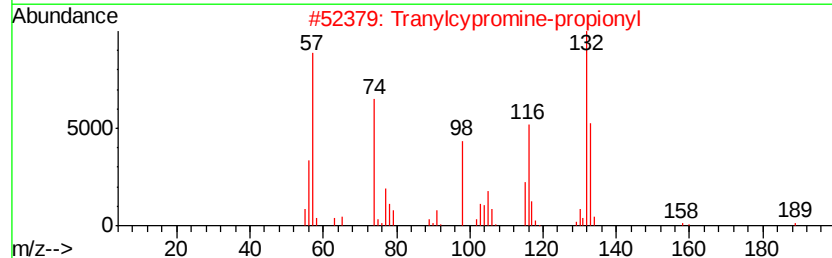
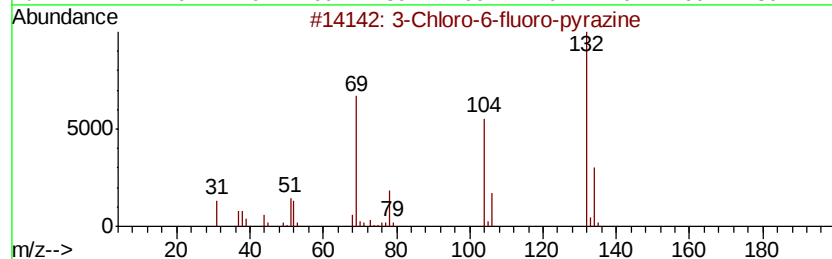
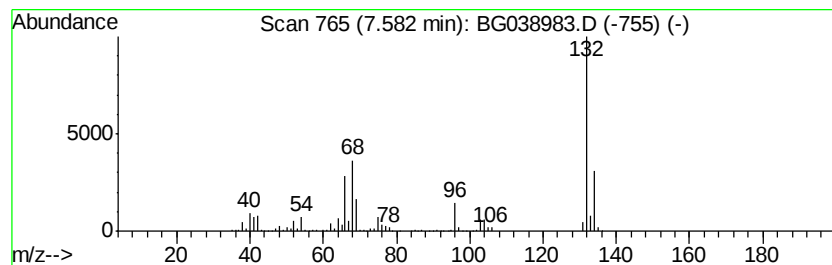
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	98.27 ng	895956	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3			5-Aminoindole	132	C8H8N2	005192-03-0	22
4			1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
5			5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.13	5.1	ng	46631	1	8.04	182354	20.0
unknown7.58	7.58	98.3	ng	895956	1	8.04	182354	20.0