

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP010820\
 Data File : BP001560.D
 Acq On : 08 Jan 2020 19:44
 Operator : JU
 Sample : PB125875BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125875BL

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_P\METHODS\8270-BP010820.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	4.834	342	348	357	rBV	136271	196275	5.61%	1.269%
2	5.293	419	426	441	rVB	720818	1043756	29.81%	6.750%
3	6.858	684	692	705	rBV	800293	1251295	35.73%	8.093%
4	7.205	740	751	762	rBV	767759	1188178	33.93%	7.685%
5	7.658	821	828	836	rVB	149629	230553	6.58%	1.491%
6	7.975	874	882	890	rBV	592652	942337	26.91%	6.095%
7	8.805	1013	1023	1034	rBV	497171	809770	23.12%	5.237%
8	10.434	1291	1300	1314	rVB	185466	303948	8.68%	1.966%
9	12.922	1716	1723	1733	rBV	1269265	1811106	51.72%	11.713%
10	14.304	1951	1958	1966	rBV2	314947	458227	13.09%	2.964%
11	15.804	2205	2213	2228	rBV	1216030	1708592	48.79%	11.050%
12	17.063	2421	2427	2436	rVB	462728	643025	18.36%	4.159%
13	19.651	2861	2867	2872	rBV	2900801	3501772	100.00%	22.648%
14	21.163	3119	3124	3129	rBV	586960	736814	21.04%	4.765%
15	23.404	3500	3505	3515	rVB	300596	636264	18.17%	4.115%

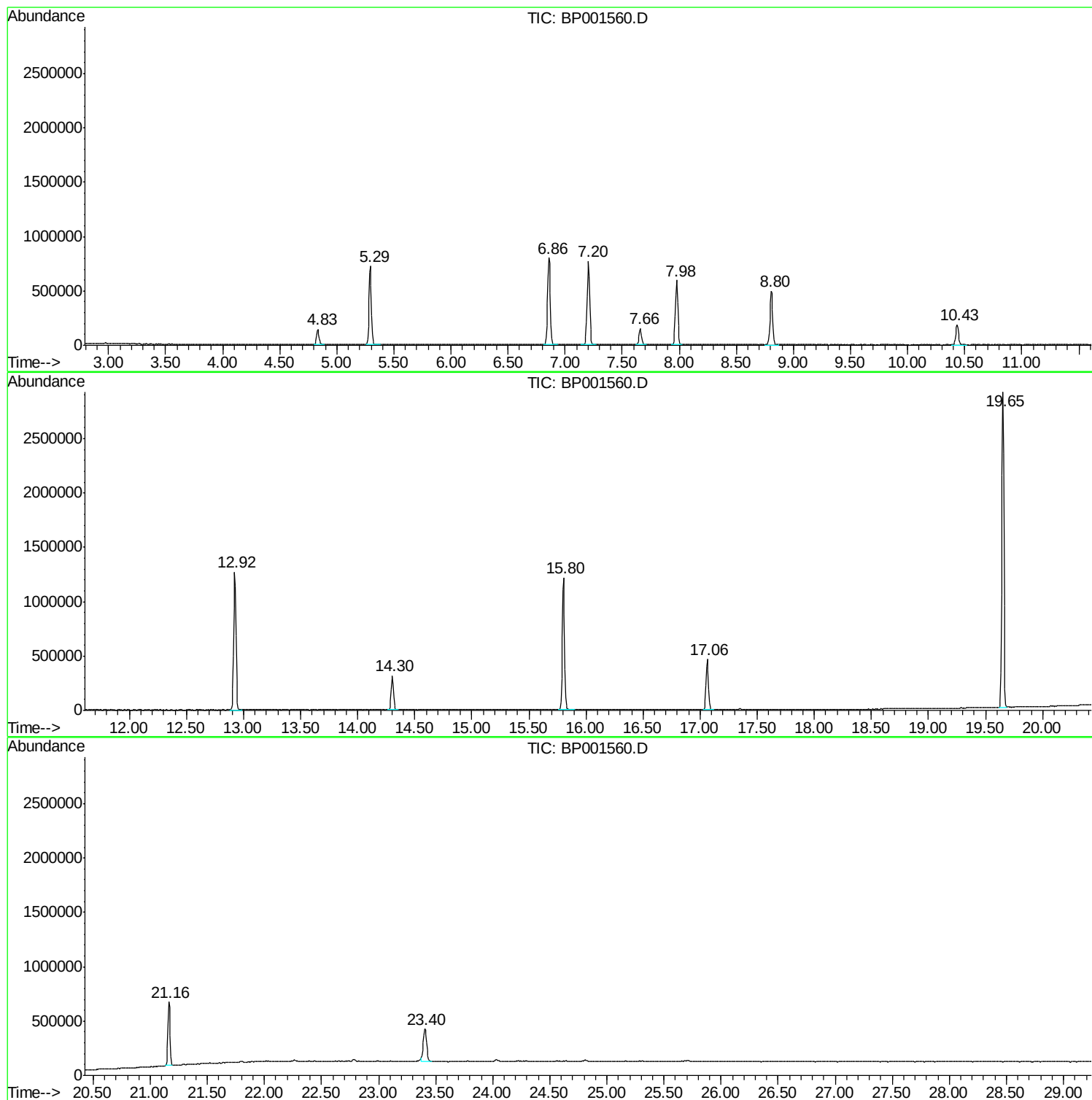
Sum of corrected areas: 15461912

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



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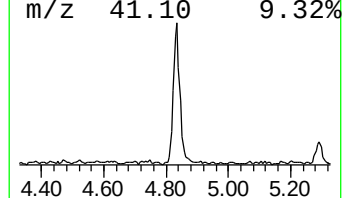
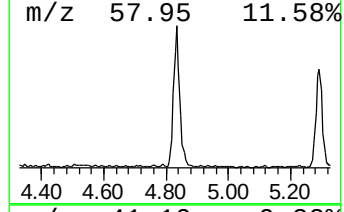
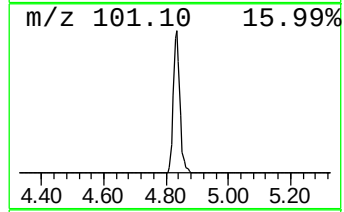
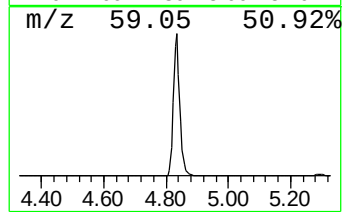
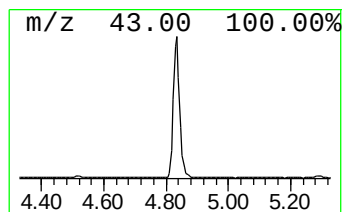
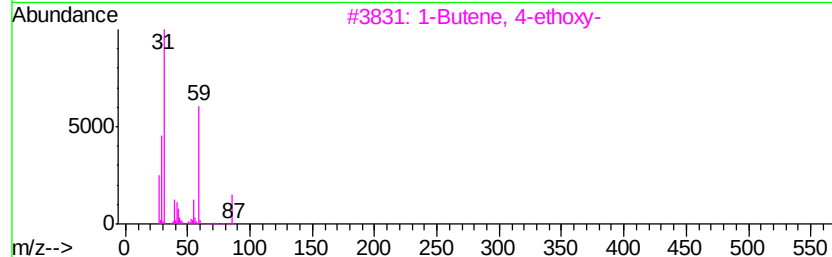
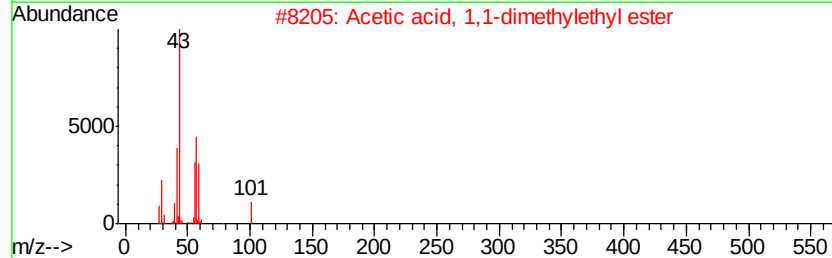
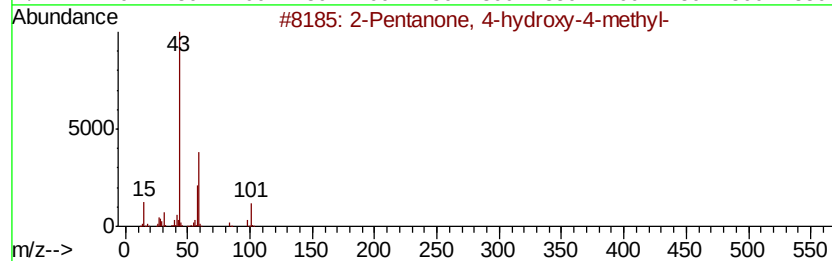
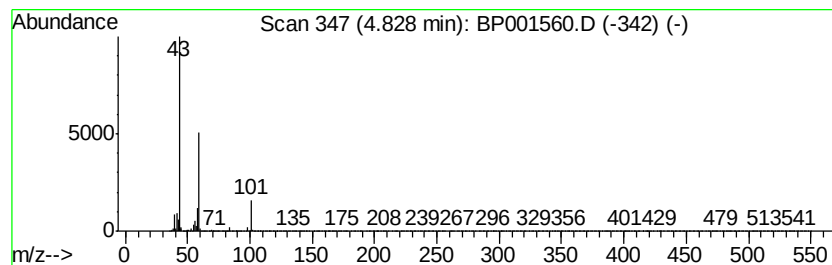
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.83	17.03 ng	196275	1,4-Dichlorobenzene-d4	7.66

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
3			1-Butene, 4-ethoxy-	100	C6H12O	044611-46-3	23
4			Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	23
5			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10



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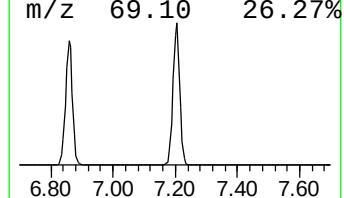
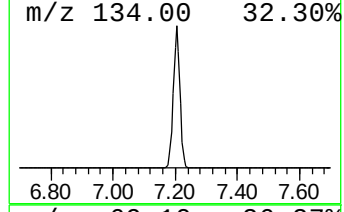
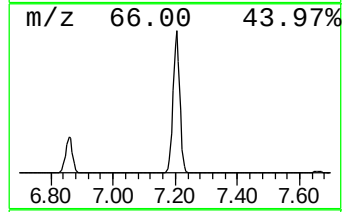
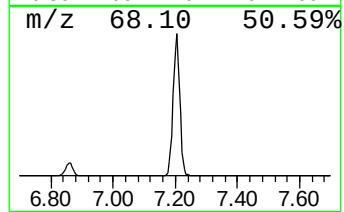
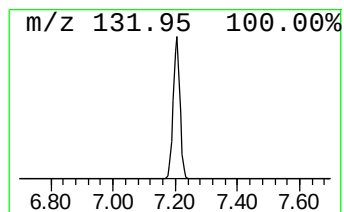
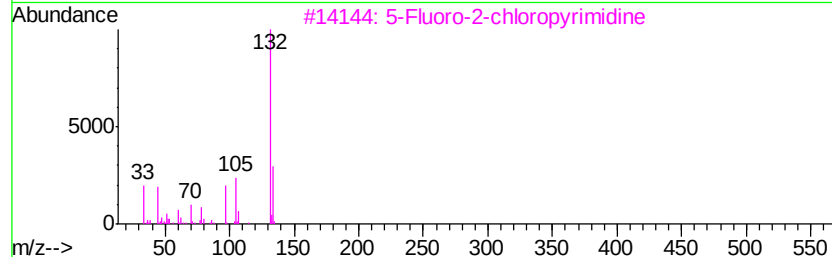
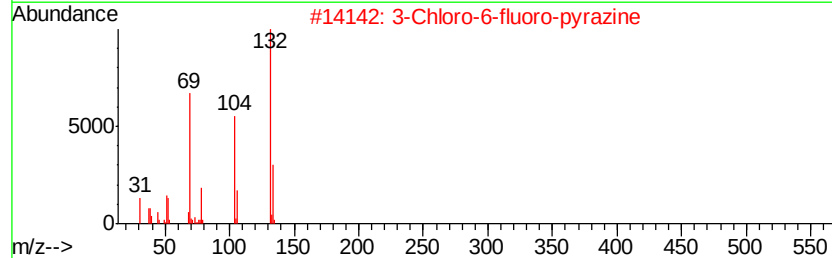
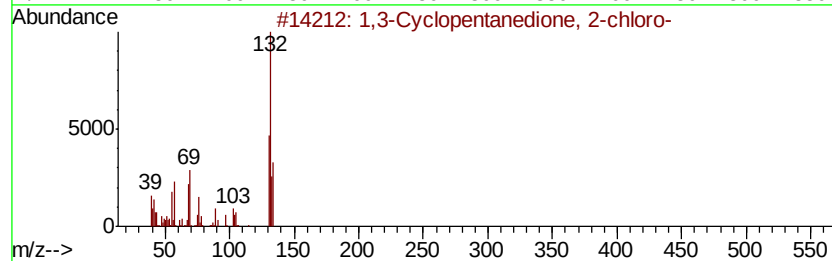
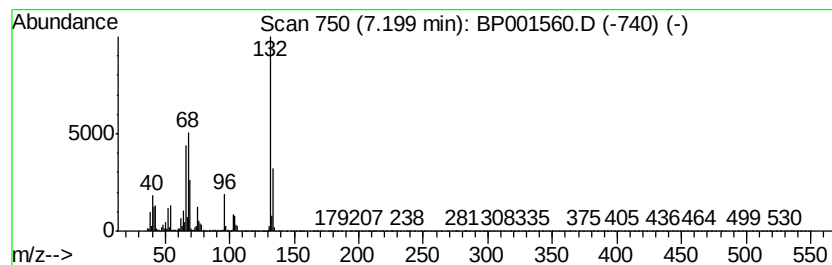
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Peak Number 2 unknown7.20 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.20	103.07 ng	1188180	1,4-Dichlorobenzene-d4	7.66

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	14
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	14
4		1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	14
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.83	17.0	ng	196275	1	7.66	230553	20.0
unknown7.20	7.20	103.1	ng	1188180	1	7.66	230553	20.0