

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN110918\
 Data File : BN003489.D
 Acq On : 10 Nov 2018 01:13
 Operator : MJ/SJ
 Sample : PB114671BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PB114671BL

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_N\METHODS\8270-BN102418.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	4.958	330	335	344	rBV	78900	110946	2.36%	0.421%
2	5.417	406	413	425	rVB	1274866	1815116	38.64%	6.892%
3	7.005	675	683	700	rBV	1220442	1892943	40.29%	7.188%
4	7.375	739	746	756	rBV	1307111	2041162	43.45%	7.751%
5	7.846	820	826	833	rBV	334909	520551	11.08%	1.977%
6	8.169	874	881	890	rVB	1501423	2357437	50.18%	8.951%
7	9.016	1017	1025	1035	rBV	1064319	1738172	37.00%	6.600%
8	10.646	1295	1302	1314	rBV	465861	770702	16.41%	2.926%
9	13.104	1713	1720	1727	rBV	3248466	4697720	100.00%	17.838%
10	14.481	1948	1954	1962	rVB2	676954	1027191	21.87%	3.900%
11	15.969	2201	2207	2216	rBV	1492133	2041776	43.46%	7.753%
12	17.228	2414	2421	2427	rBV	773862	1064429	22.66%	4.042%
13	19.845	2860	2866	2871	rBV	3278078	4168542	88.74%	15.828%
14	21.404	3126	3131	3136	rBV	850059	1013188	21.57%	3.847%
15	23.716	3517	3524	3535	rVB2	553487	1075959	22.90%	4.086%

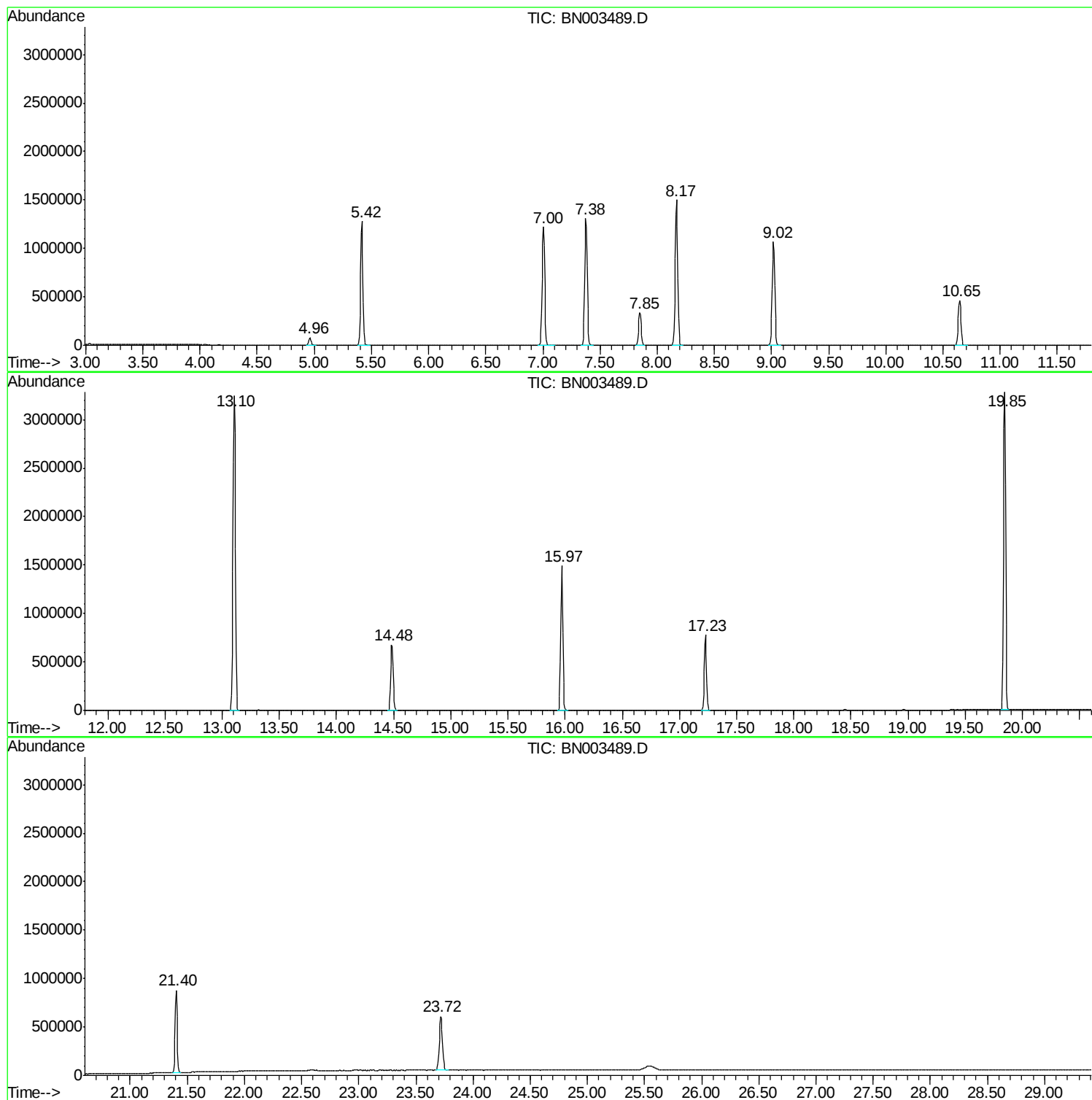
Sum of corrected areas: 26335834

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN102418.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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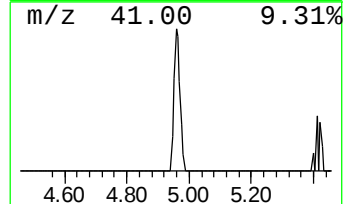
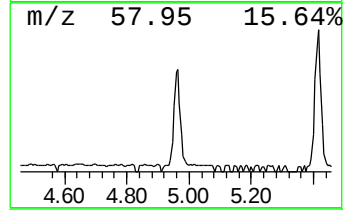
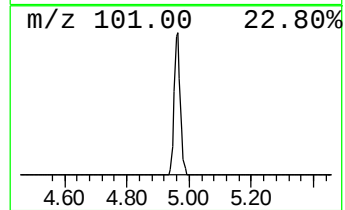
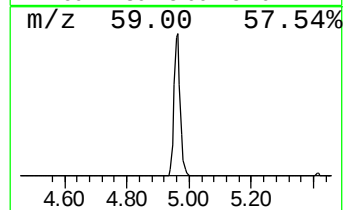
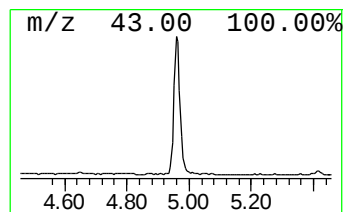
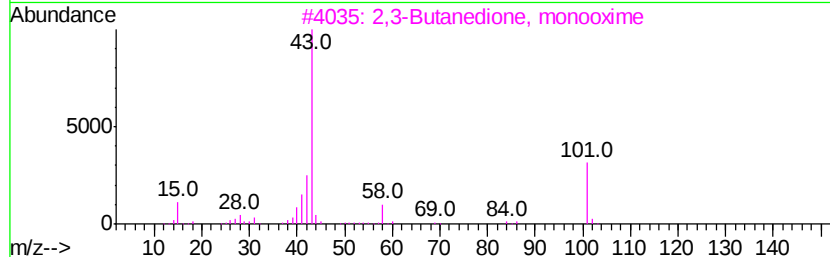
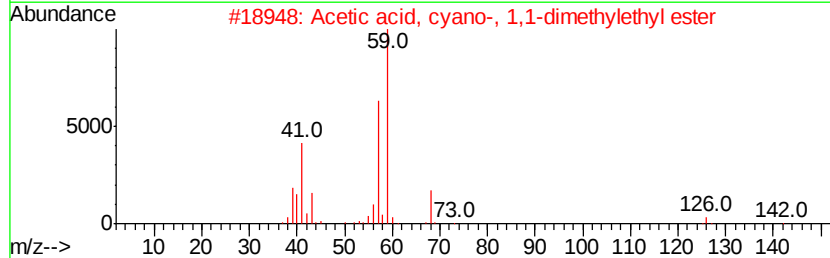
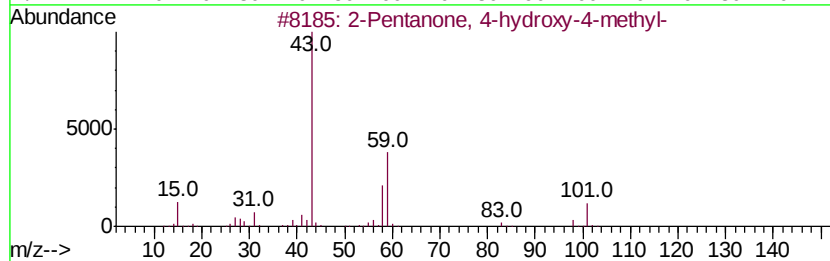
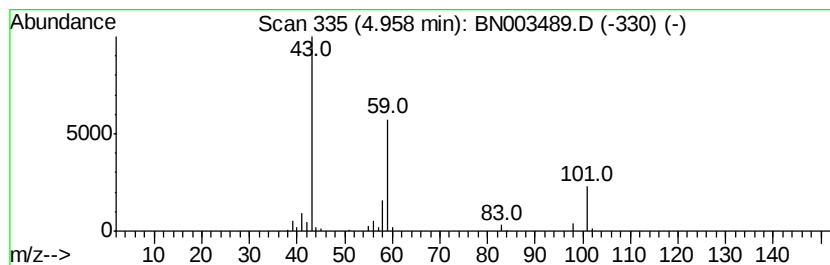
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.96	4.26 ng	110946	1,4-Dichlorobenzene-d4	7.85

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
3			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5			(+)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9



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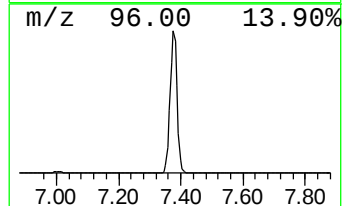
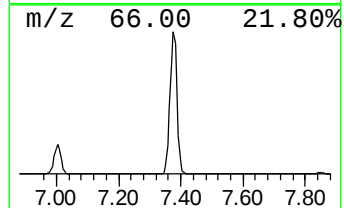
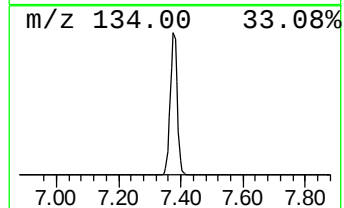
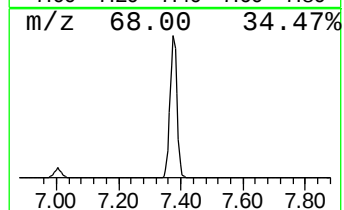
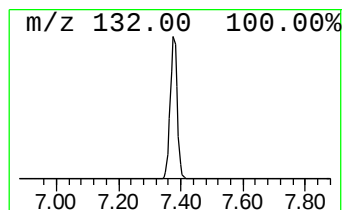
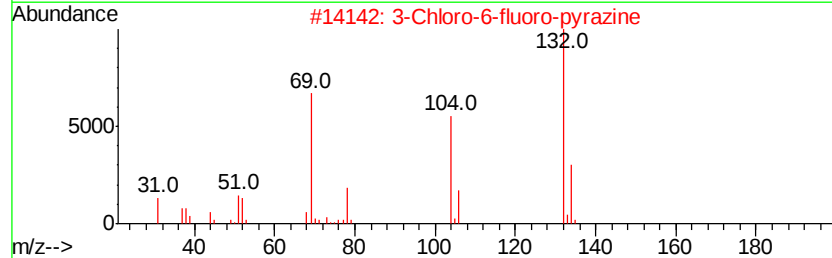
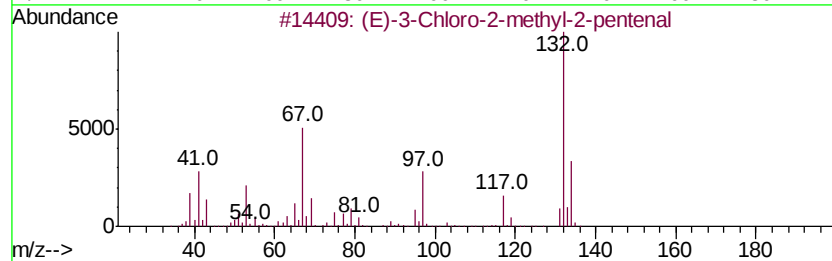
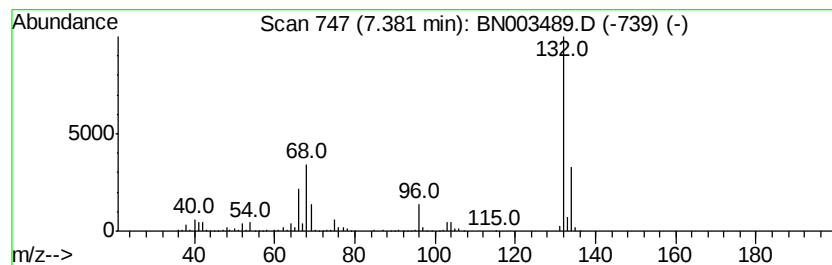
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Peak Number 2 unknown7.38 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.38	78.42 ng	2041160	1,4-Dichlorobenzene-d4	7.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	35
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
3		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		5-Aminoindole	132	C8H8N2	005192-03-0	12
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.96	4.3	ng	110946	1	7.85	520551	20.0
unknown7.38	7.38	78.4	ng	2041160	1	7.85	520551	20.0