

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF070218\
 Data File : BF107039.D
 Acq On : 2 Jul 2018 12:50
 Operator : JU/SJ
 Sample : J3761-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 256-CONCORD-COMP

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_F\METHODS\8270-BF061918.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	5.190	481	485	495	rBV	110989	130438	10.43%	1.187%
2	5.601	551	555	568	rBV	1197529	1118296	89.41%	10.176%
3	6.601	721	725	736	rBV	1090146	1131687	90.49%	10.297%
4	6.737	744	748	751	rBV	1306091	1162010	92.91%	10.573%
5	6.954	782	785	789	rBV	420877	338656	27.08%	3.081%
6	7.113	808	812	815	rBV	1108966	913840	73.07%	8.315%
7	7.519	877	881	895	rBV	724594	709143	56.70%	6.453%
8	8.236	1000	1003	1019	rBV	410624	422862	33.81%	3.848%
9	9.313	1182	1186	1201	rVB	1417177	1250684	100.00%	11.380%
10	9.707	1249	1253	1261	rVB	98900	94256	7.54%	0.858%
11	10.001	1299	1303	1315	rVB	467455	445946	35.66%	4.058%
12	10.795	1434	1438	1445	rBV	806736	737591	58.98%	6.711%
13	10.877	1449	1452	1458	rVB	16003	17338	1.39%	0.158%
14	11.495	1553	1557	1564	rBV	485360	423264	33.84%	3.851%
15	12.948	1801	1804	1807	rVB	16424	16216	1.30%	0.148%
16	13.077	1822	1826	1829	rBV	1402931	1199857	95.94%	10.918%
17	13.954	1972	1975	1978	rBV	44397	50089	4.00%	0.456%
18	14.154	2005	2009	2014	rBV	497844	507340	40.57%	4.616%
19	15.707	2268	2273	2279	rVB	220656	320506	25.63%	2.916%

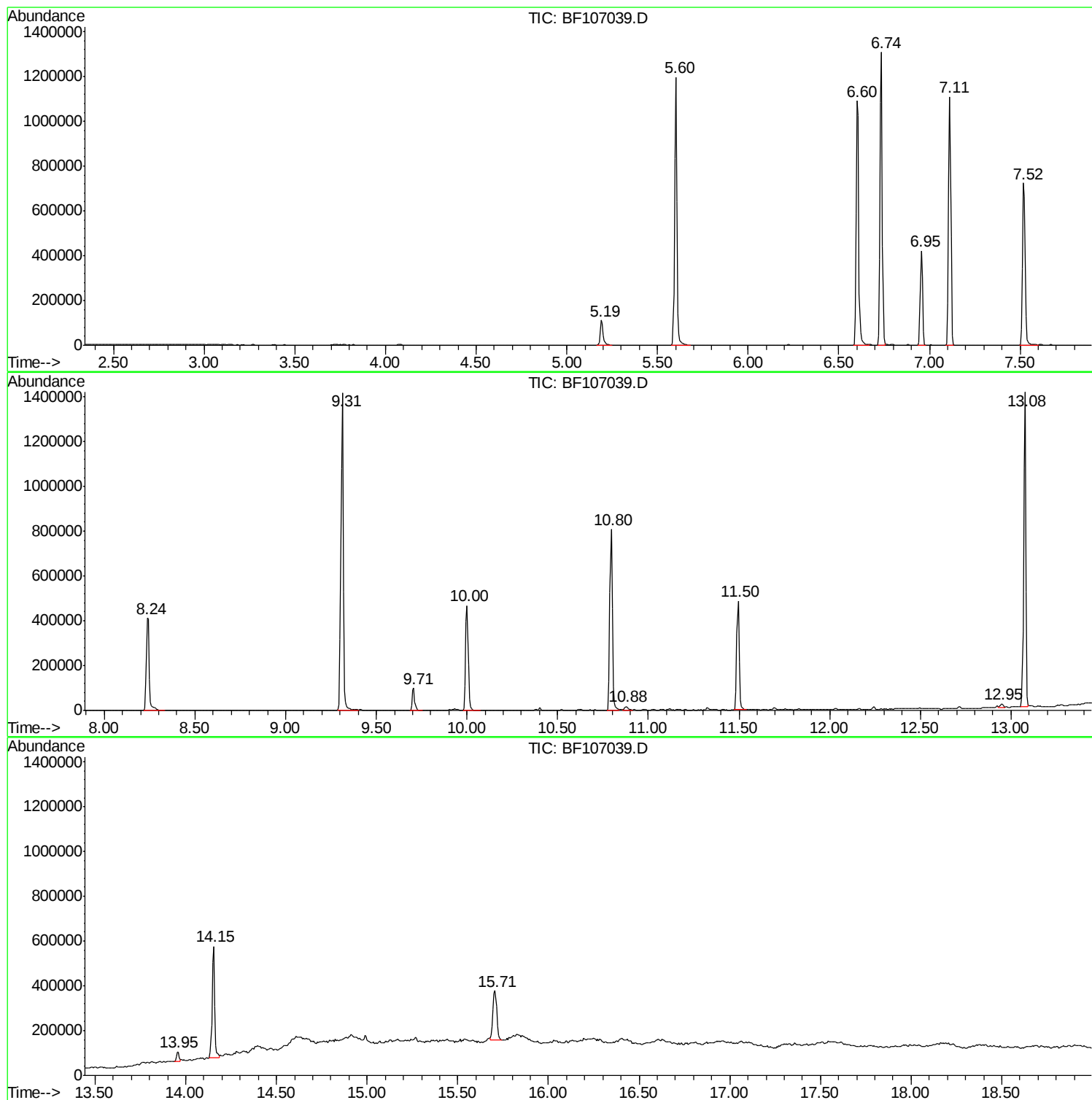
Sum of corrected areas: 10990019

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.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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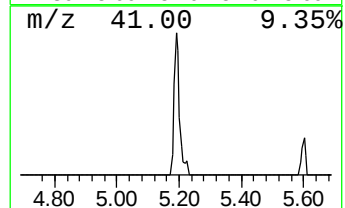
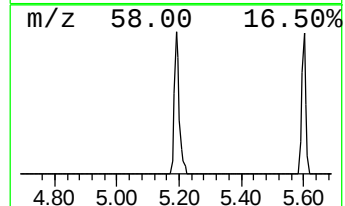
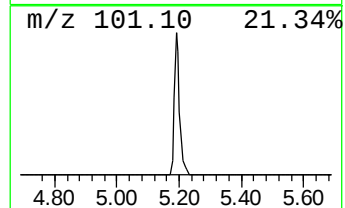
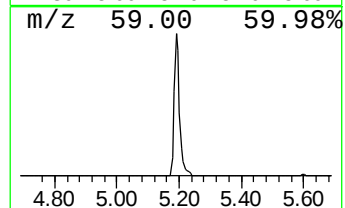
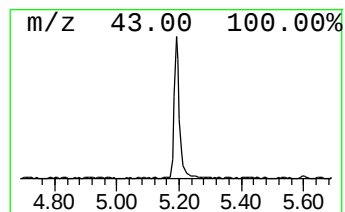
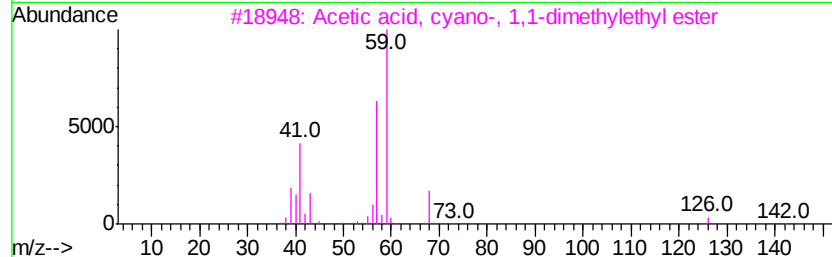
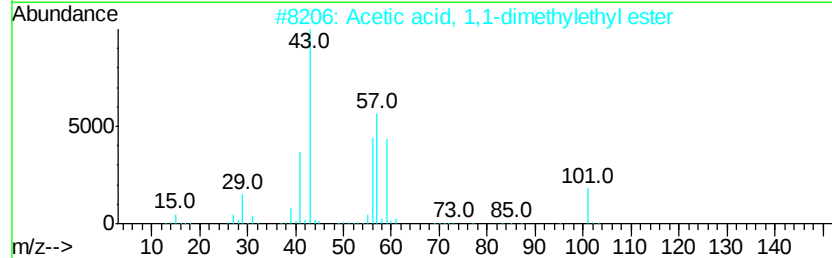
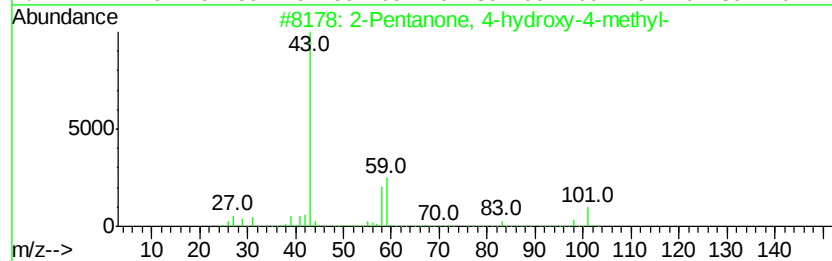
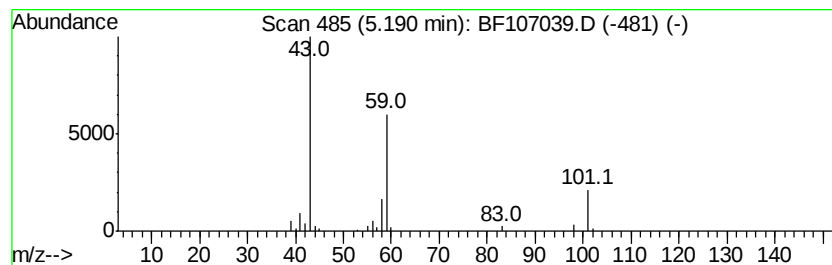
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.19	7.70 ng	130438	1,4-Dichlorobenzene-d4	6.95

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	39
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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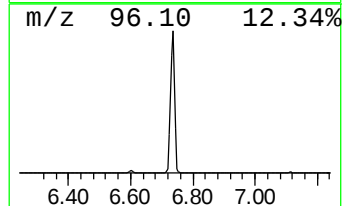
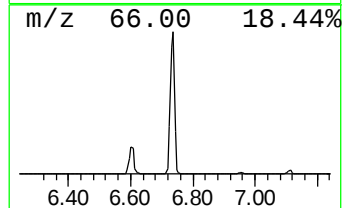
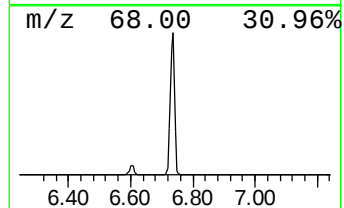
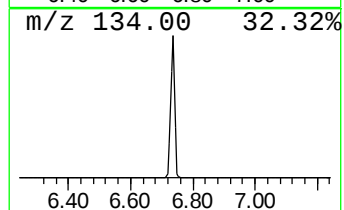
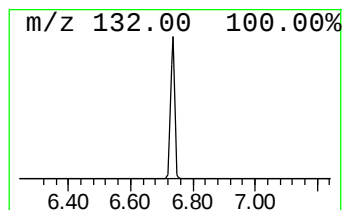
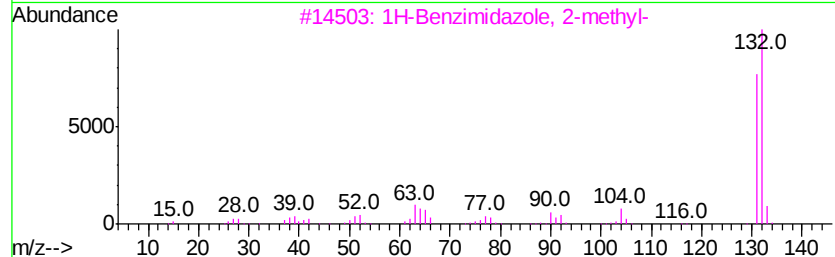
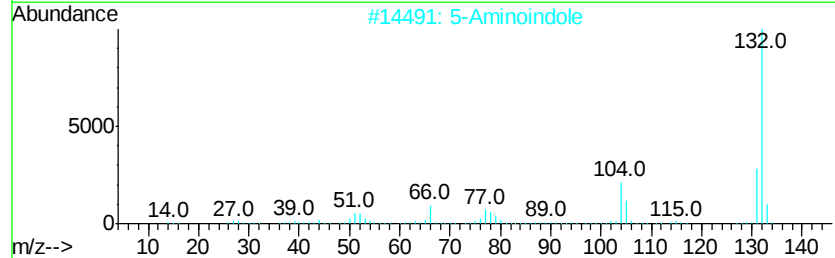
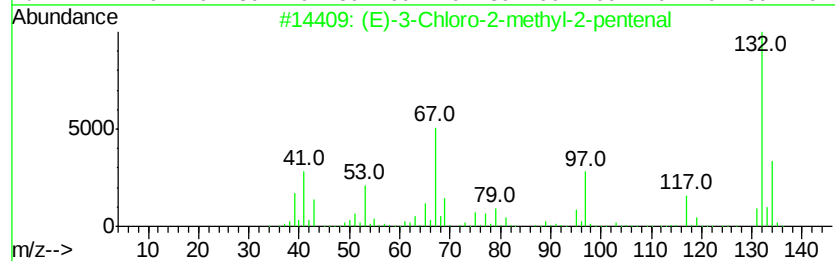
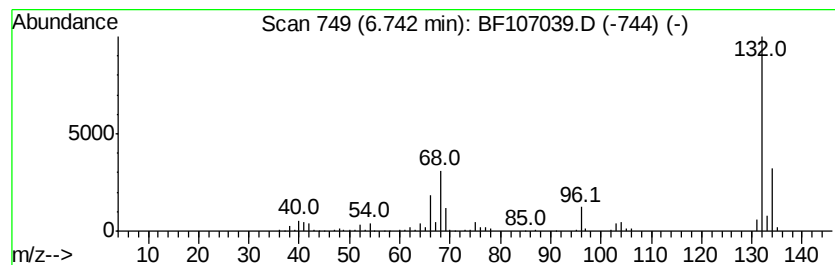
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	68.62 ng	1162010	1,4-Dichlorobenzene-d4	6.95

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Xenon	132	Xe	007440-63-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.19	7.7	ng	130438	1	6.95	338656	20.0
unknown6.74	6.74	68.6	ng	1162010	1	6.95	338656	20.0