

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061518\
 Data File : BF106522.D
 Acq On : 16 Jun 2018 5:37
 Operator : JU/SJ
 Sample : J3447-04
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-32

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-BF061118.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	5.195	482	486	494	rVB	75162	82503	3.12%	0.451%
2	5.607	553	556	567	rBV	1302782	1338713	50.54%	7.323%
3	6.607	722	726	737	rBV	978001	894857	33.79%	4.895%
4	6.742	745	749	752	rBV	2333515	2078921	78.49%	11.372%
5	6.966	783	787	790	rBV	815716	644039	24.32%	3.523%
6	7.125	809	814	817	rBV	2178973	1992007	75.21%	10.897%
7	7.531	878	883	886	rBV	1641641	1590169	60.04%	8.699%
8	8.248	1001	1005	1008	rBV	883124	698616	26.38%	3.822%
9	9.319	1183	1187	1191	rBV	2491826	2436165	91.98%	13.326%
10	10.007	1300	1304	1310	rBV	711475	628746	23.74%	3.439%
11	10.801	1434	1439	1442	rBV	1413016	1354375	51.14%	7.409%
12	11.495	1554	1557	1561	rBV	708108	622465	23.50%	3.405%
13	13.083	1823	1827	1830	rBV	2774207	2648561	100.00%	14.488%
14	14.148	2004	2008	2013	rBV	851017	736192	27.80%	4.027%
15	15.001	2150	2153	2156	rBV	28824	30626	1.16%	0.168%
16	15.683	2263	2269	2277	rBV	370580	503914	19.03%	2.757%

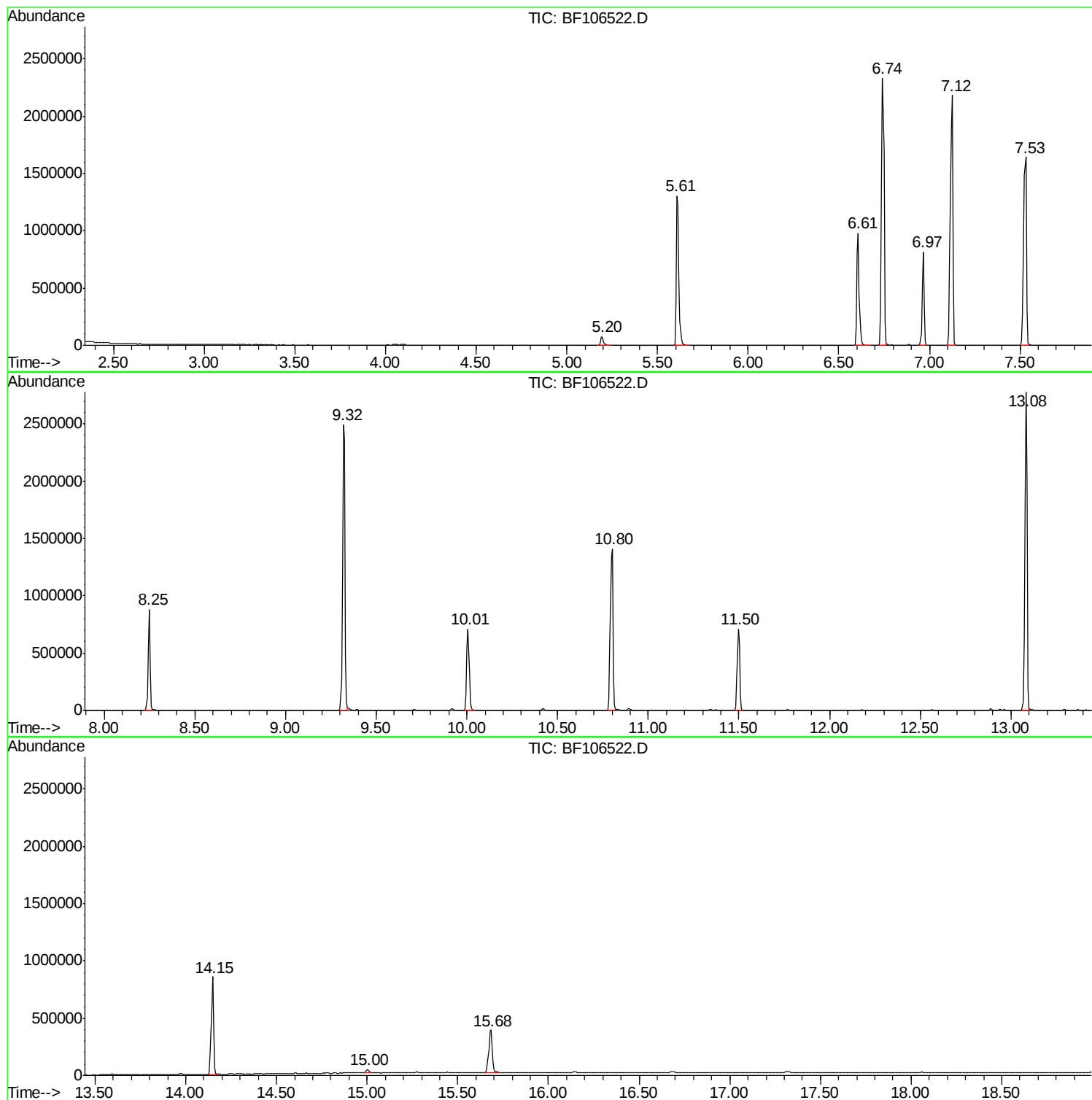
Sum of corrected areas: 18280869

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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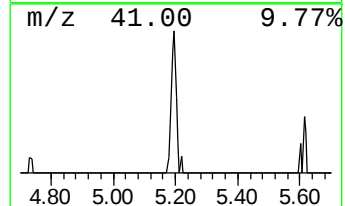
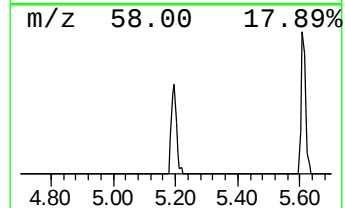
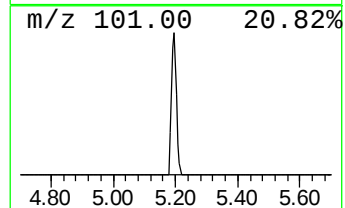
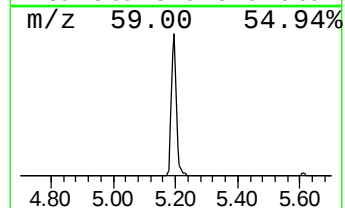
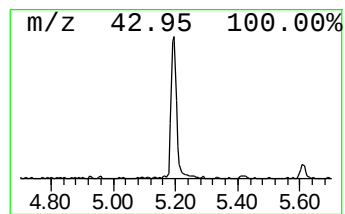
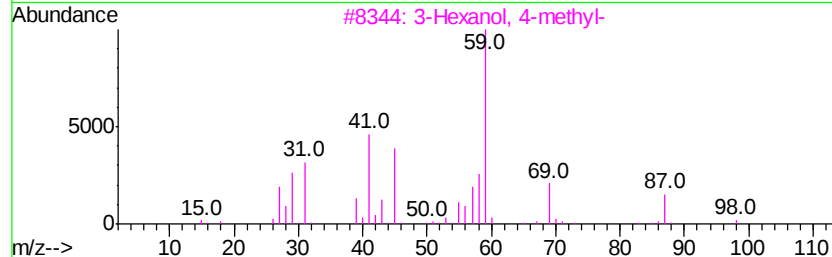
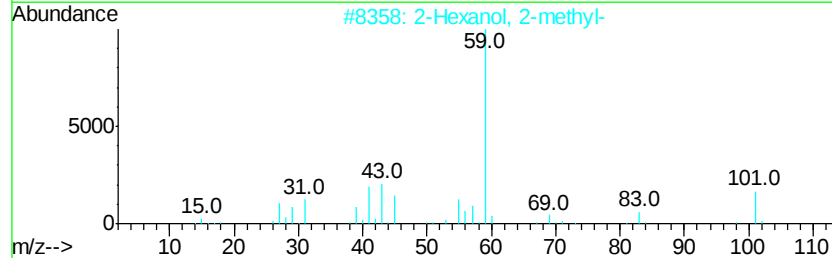
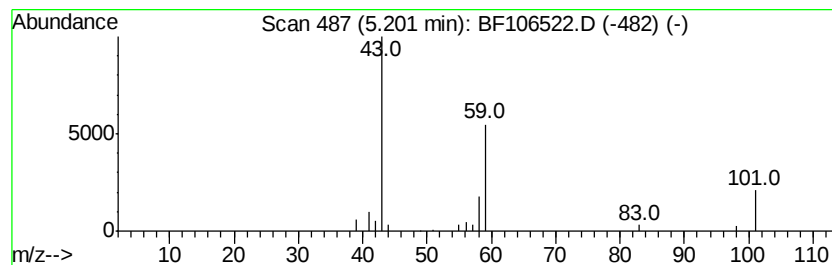
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	2.56 ng	82503	1,4-Dichlorobenzene-d4	6.97

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			2-Hexanol, 2-methyl-	116	C7H16O	000625-23-0	28
3			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
4			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
5			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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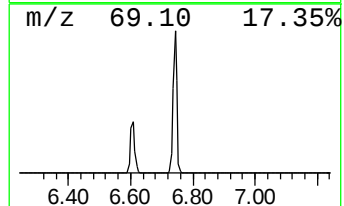
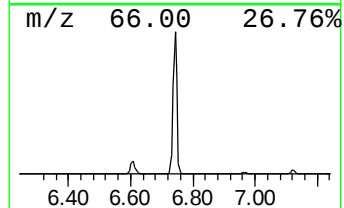
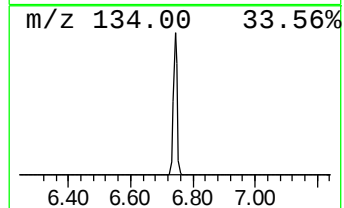
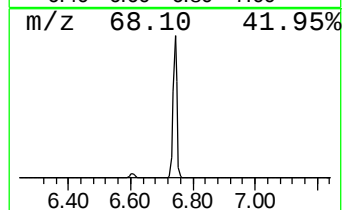
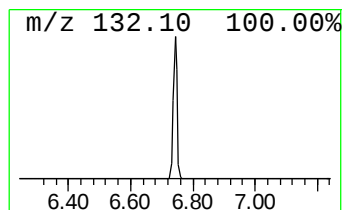
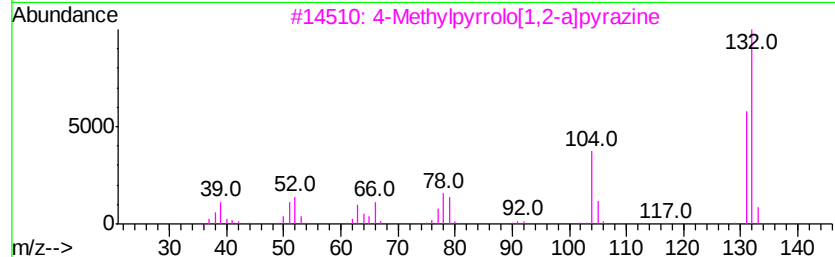
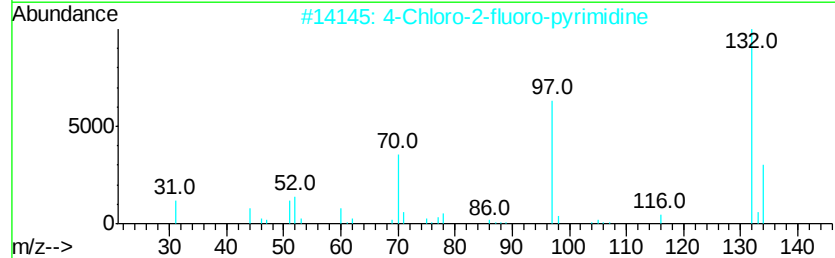
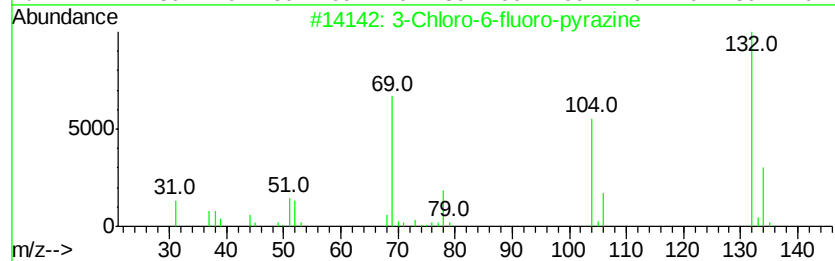
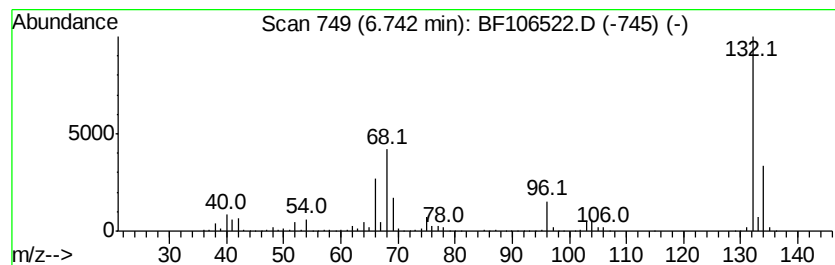
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Peak Number 2 unknown6.74 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	64.56 ng	2078920	1,4-Dichlorobenzene-d4	6.97

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
3		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	2.6	ng	82503	1	6.97	644039	20.0
unknown6.74	6.74	64.6	ng	2078920	1	6.97	644039	20.0