

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062124\
 Data File : BF138279.D
 Acq On : 21 Jun 2024 11:50
 Operator : CG\JU
 Sample : PB161600BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB161600BL

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-BF061224.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138279.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.252	17	28	33	rVB	137368	206433	1.89%	0.328%
2	3.563	244	251	266	rBV2	52303	142038	1.30%	0.226%
3	5.128	510	517	530	rVB	319393	460491	4.22%	0.732%
4	5.516	577	583	586	rBV	6335218	7313517	66.99%	11.625%
5	6.516	746	753	756	rBV	6028133	7262707	66.52%	11.544%
6	6.881	812	815	819	rBV	2348756	1893783	17.35%	3.010%
7	7.445	905	911	914	rBV	4465649	4956819	45.40%	7.879%
8	8.163	1028	1033	1036	rBV	3012158	2584340	23.67%	4.108%
9	8.228	1039	1044	1052	rVB2	93634	109617	1.00%	0.174%
10	9.245	1210	1217	1220	rBV	8667631	9413375	86.22%	14.963%
11	9.916	1326	1331	1335	rBV	3562573	3139638	28.76%	4.991%
12	10.710	1461	1466	1469	rBV	6449042	6178135	56.59%	9.820%
13	11.404	1580	1584	1587	rBV	4162369	3377184	30.93%	5.368%
14	12.998	1849	1855	1858	rBV	9660420	10917544	100.00%	17.354%
15	14.039	2028	2032	2043	rVB	3007735	2667852	24.44%	4.241%
16	15.510	2277	2282	2298	rBV	1503018	2151519	19.71%	3.420%
17	17.121	2549	2556	2559	rBV3	53174	136650	1.25%	0.217%

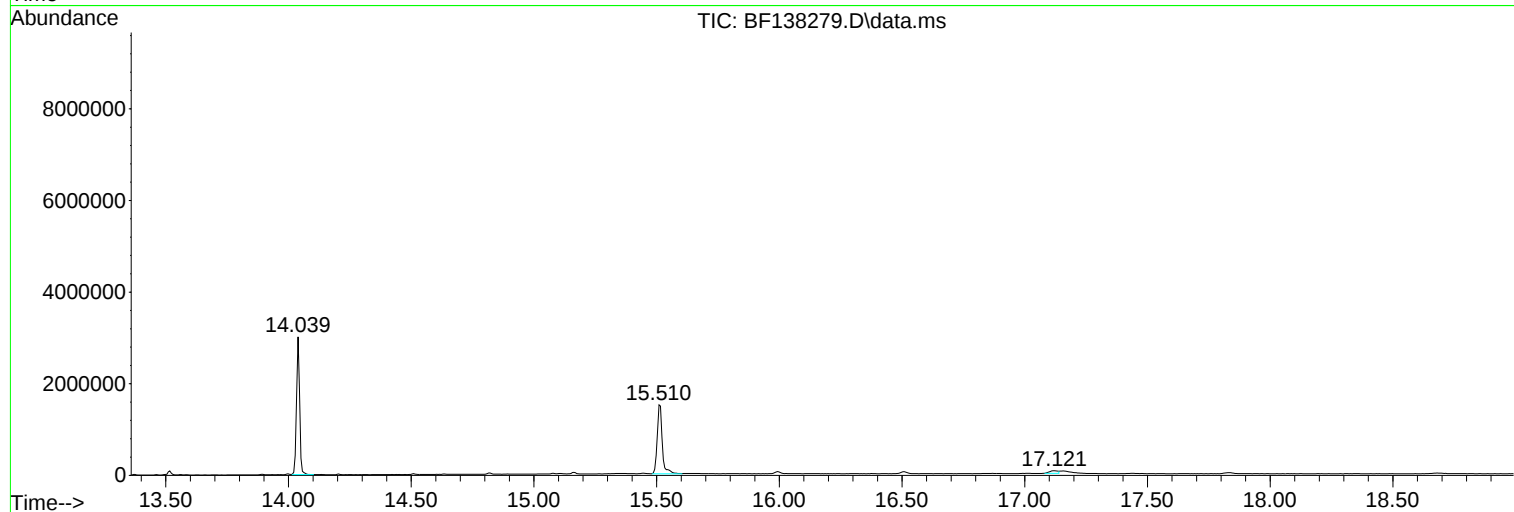
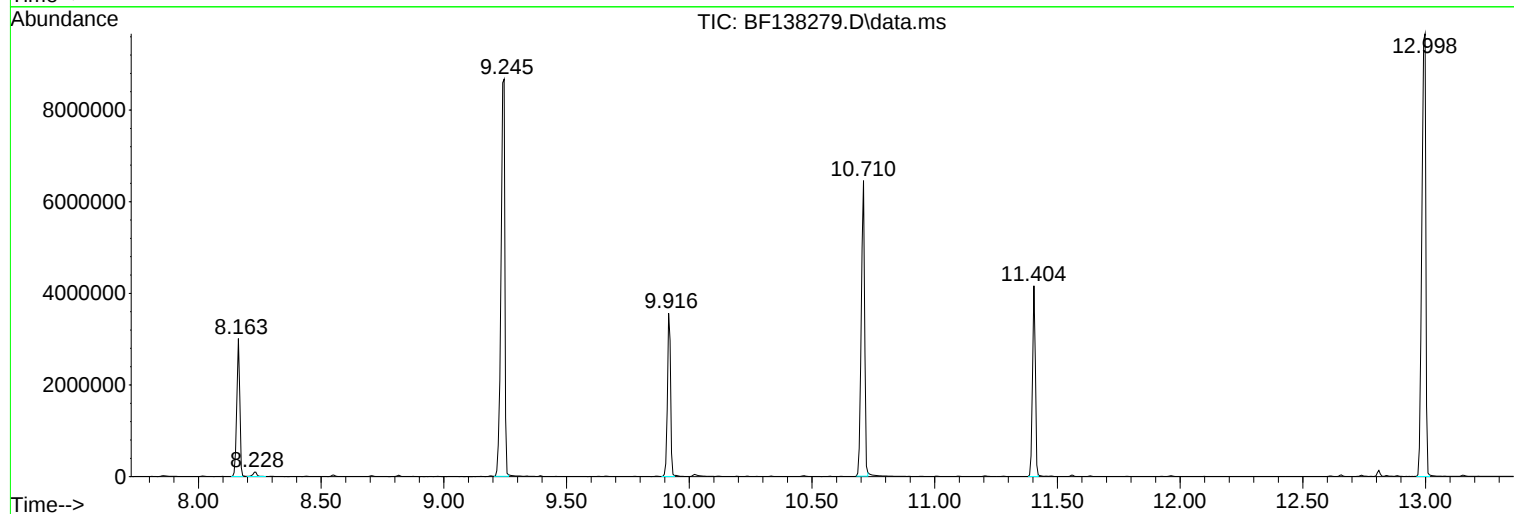
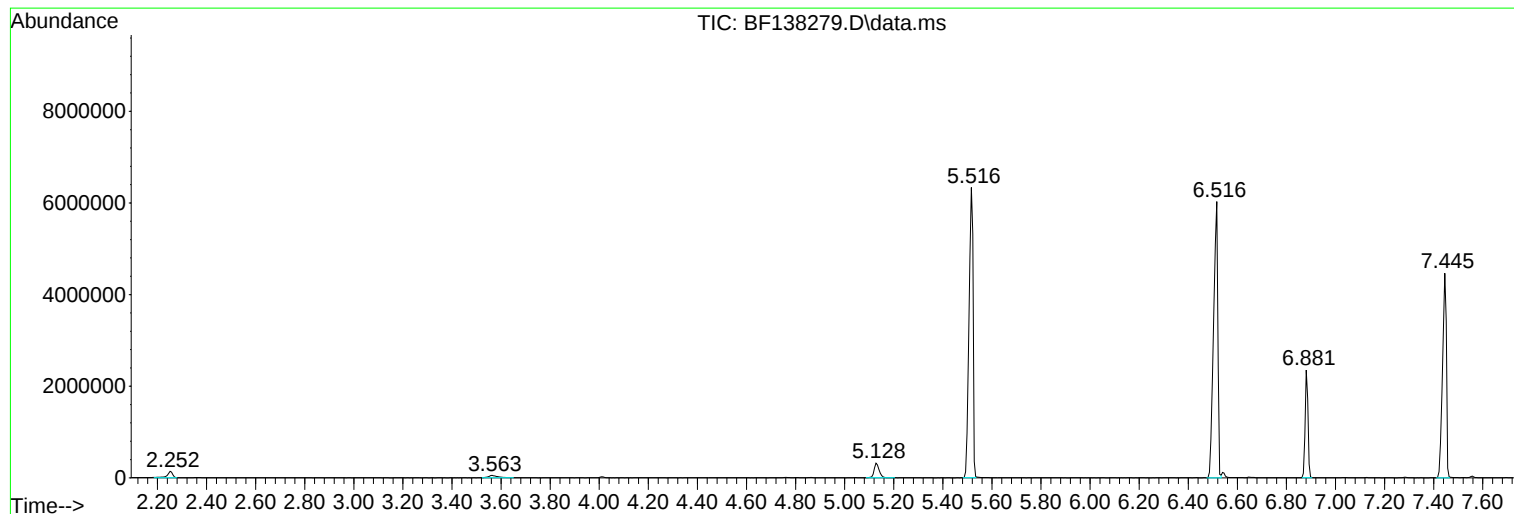
Sum of corrected areas: 62911642

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061224.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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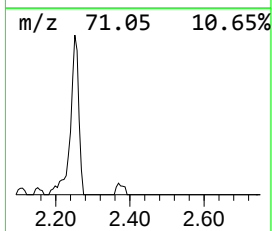
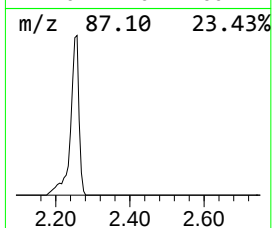
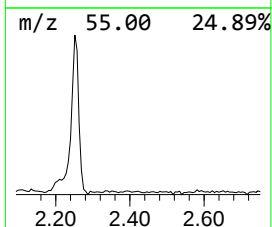
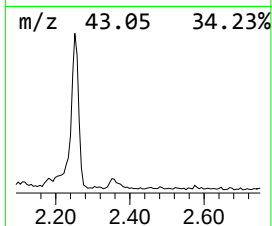
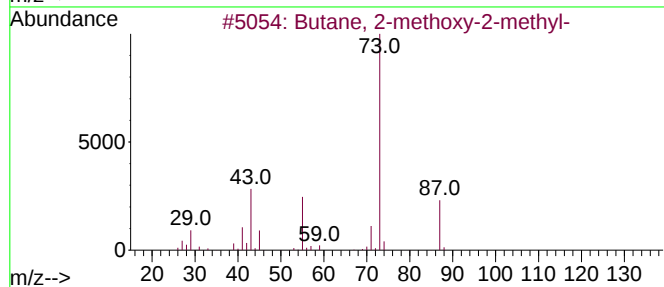
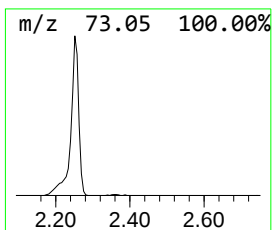
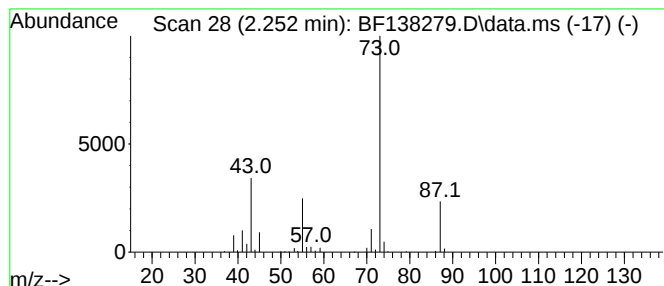
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.252	2.18 ng	206433	1,4-Dichlorobenzene-d4	6.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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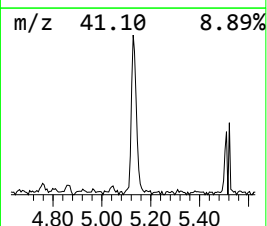
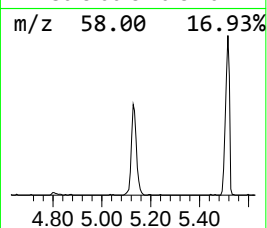
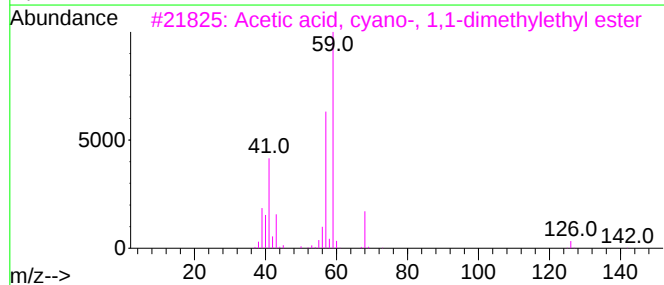
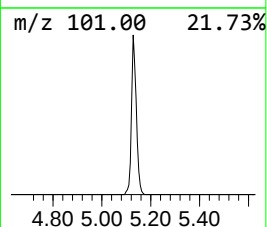
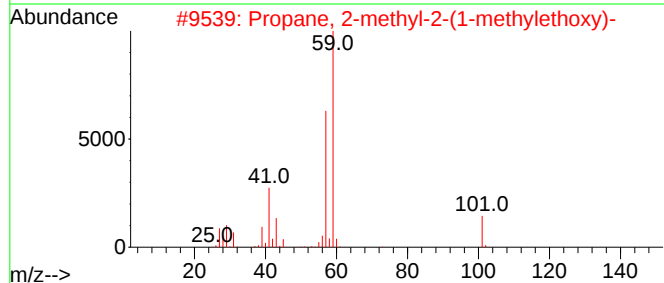
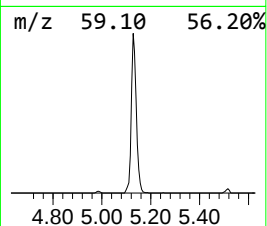
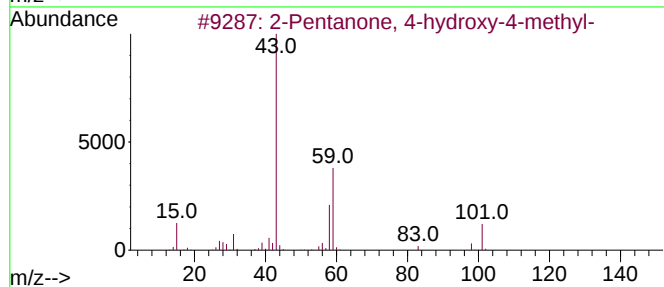
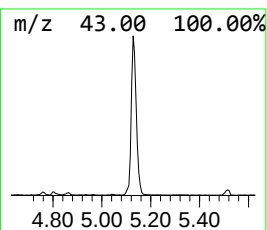
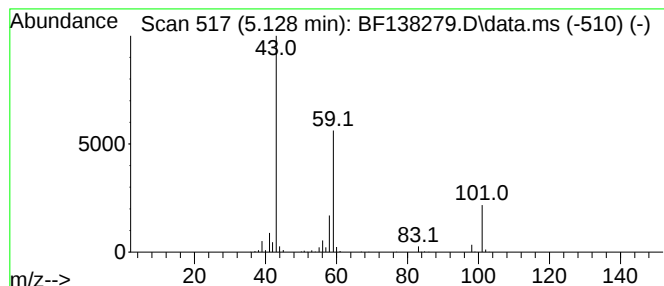
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.128	4.86 ng	460491	1,4-Dichlorobenzene-d4	6.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	45		
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	23		
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17		
4	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	10		
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9		



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.252	2.2	ng	206433	1	6.881	1893780	20.0
2-Pentanone, 4-...	5.128	4.9	ng	460491	1	6.881	1893780	20.0