

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071423\
 Data File : BF134284.D
 Acq On : 14 Jul 2023 11:22
 Operator : CG\JU
 Sample : PB154028BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB154028BL

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-BF062623.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF134284.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.146	4	10	19	rVB	56183	84231	2.38%	0.429%
2	5.099	507	512	521	rBV	45860	67632	1.91%	0.345%
3	5.481	573	577	592	rBV	2381781	2473716	69.79%	12.601%
4	6.487	743	748	766	rBV	2367600	2468194	69.63%	12.573%
5	6.840	805	808	820	rBV	605151	506107	14.28%	2.578%
6	7.410	900	905	908	rBV	1746875	1586879	44.77%	8.084%
7	8.122	1022	1026	1032	rBV	815746	641221	18.09%	3.266%
8	9.204	1205	1210	1213	rBV	3288452	3025562	85.35%	15.412%
9	9.881	1321	1325	1338	rBV	915379	735108	20.74%	3.745%
10	10.675	1456	1460	1479	rBV	2185534	2171988	61.27%	11.064%
11	11.375	1575	1579	1592	rBV	933151	821185	23.17%	4.183%
12	12.980	1847	1852	1855	rBV	3631547	3544735	100.00%	18.057%
13	14.039	2028	2032	2044	rBV	843312	764915	21.58%	3.897%
14	15.545	2283	2288	2298	rBV	550269	739262	20.86%	3.766%

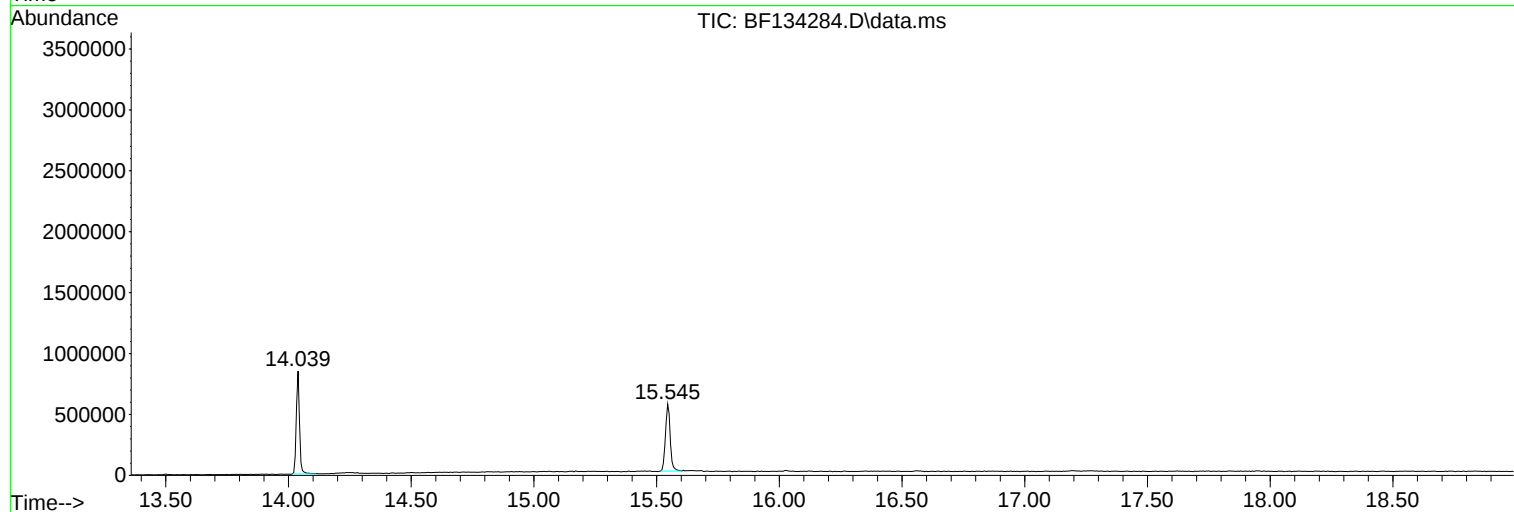
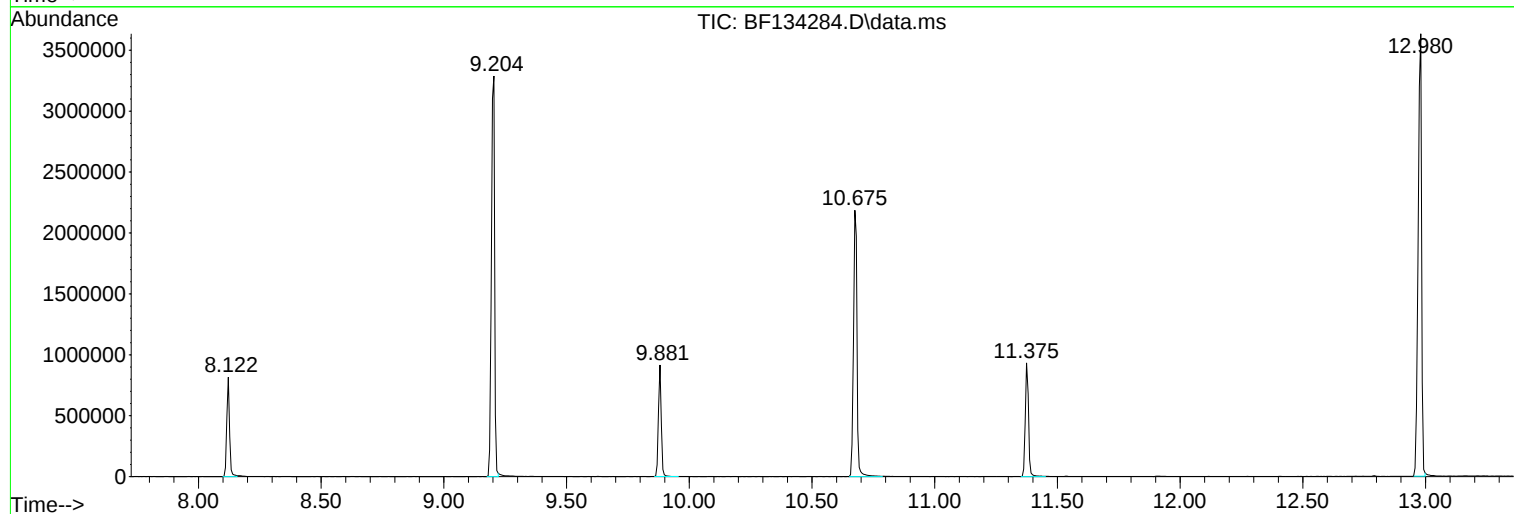
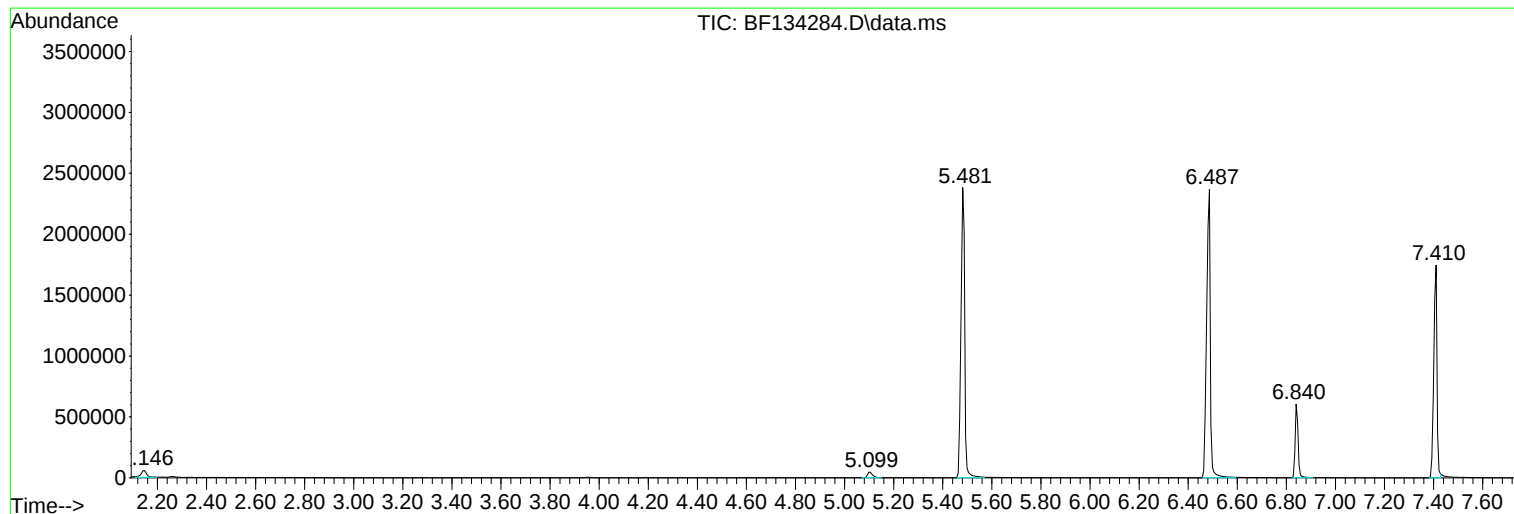
Sum of corrected areas: 19630735

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062623.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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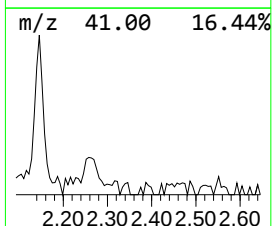
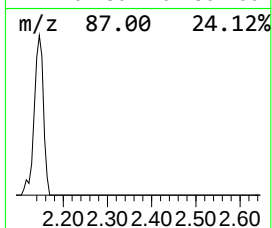
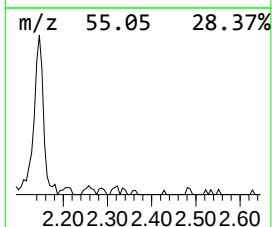
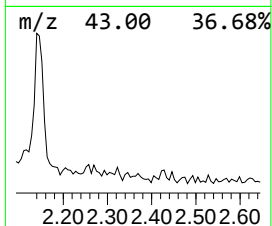
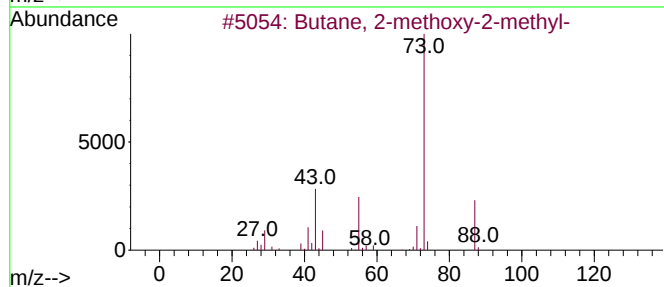
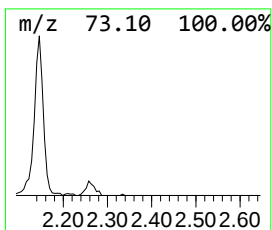
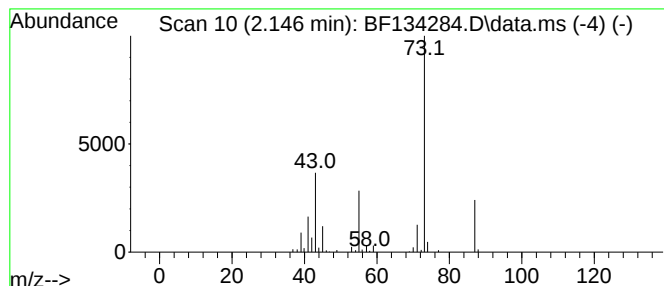
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.146	3.33 ng	84231	1,4-Dichlorobenzene-d4	6.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
4			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	37
5			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37



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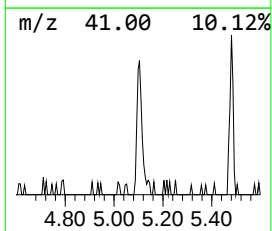
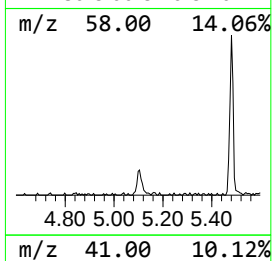
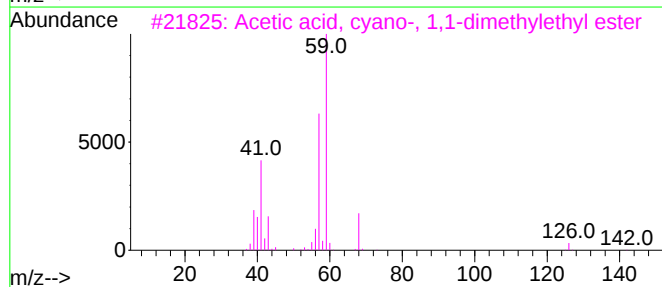
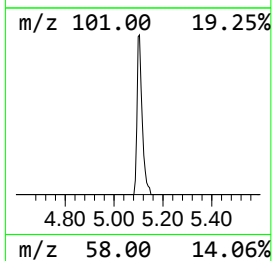
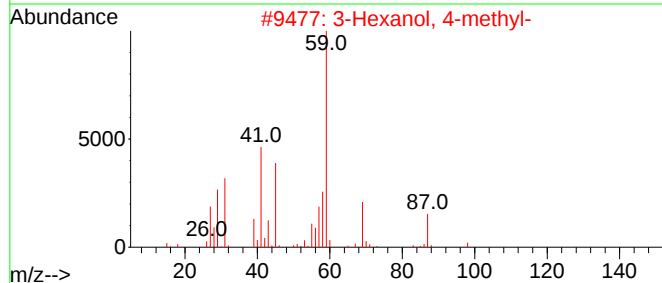
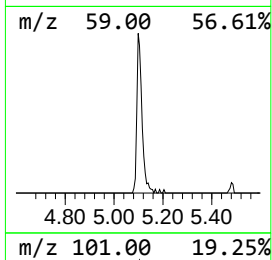
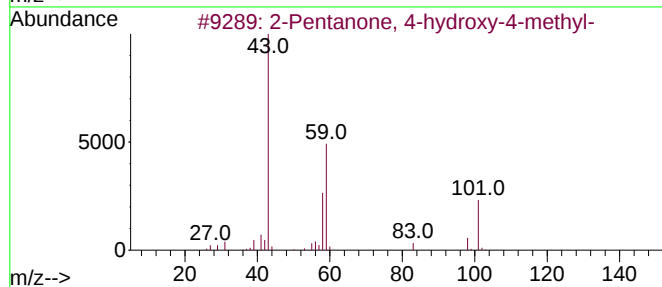
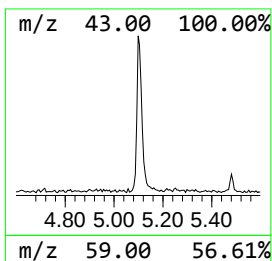
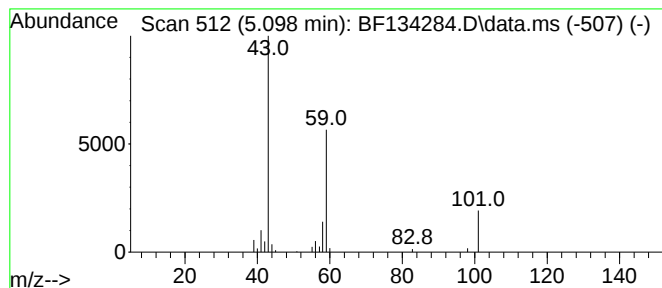
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.098	2.67 ng	67632	1,4-Dichlorobenzene-d4			6.840
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	56
2	3-Hexanol, 4-methyl-		116	C7H16O	000615-29-2	25
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	2,3-Butanedione, monooxime		101	C4H7NO2	000057-71-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.146	3.3	ng	84231	1	6.840	506107	20.0
2-Pentanone, 4-...	5.098	2.7	ng	67632	1	6.840	506107	20.0