

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072424\
 Data File : BF138639.D
 Acq On : 24 Jul 2024 14:34
 Operator : RC/JU
 Sample : PB162256BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB162256BL

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-BF070924.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138639.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.216	16	21	31	rVB	45510	77946	2.38%	0.397%
2	5.098	505	511	524	rVB	119503	200773	6.14%	1.022%
3	5.487	571	577	583	rBV	1995362	2737737	83.67%	13.941%
4	6.486	741	747	753	rBV	1873015	2793636	85.38%	14.225%
5	6.851	804	809	814	rVB	404052	486354	14.86%	2.477%
6	7.416	898	905	910	rBV	1306810	1875354	57.32%	9.549%
7	8.133	1020	1027	1032	rBV	489424	660220	20.18%	3.362%
8	9.210	1202	1210	1215	rBV	2242221	3188376	97.45%	16.236%
9	9.886	1319	1325	1339	rBV	579620	742940	22.71%	3.783%
10	10.674	1453	1459	1488	rBV	1403120	1905853	58.25%	9.705%
11	11.369	1572	1577	1584	rBV	625135	769778	23.53%	3.920%
12	12.957	1841	1847	1852	rBV	2439748	3271936	100.00%	16.661%
13	14.004	2020	2025	2036	rBV	414415	532874	16.29%	2.713%
14	15.468	2267	2274	2295	rBV	235760	394494	12.06%	2.009%

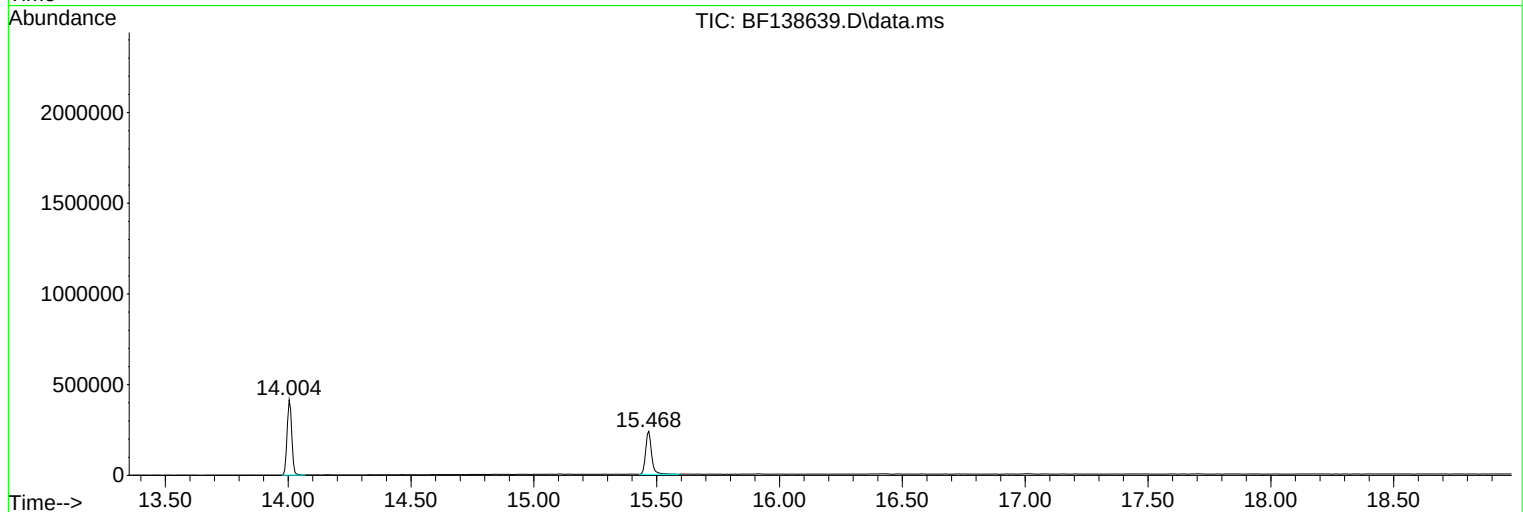
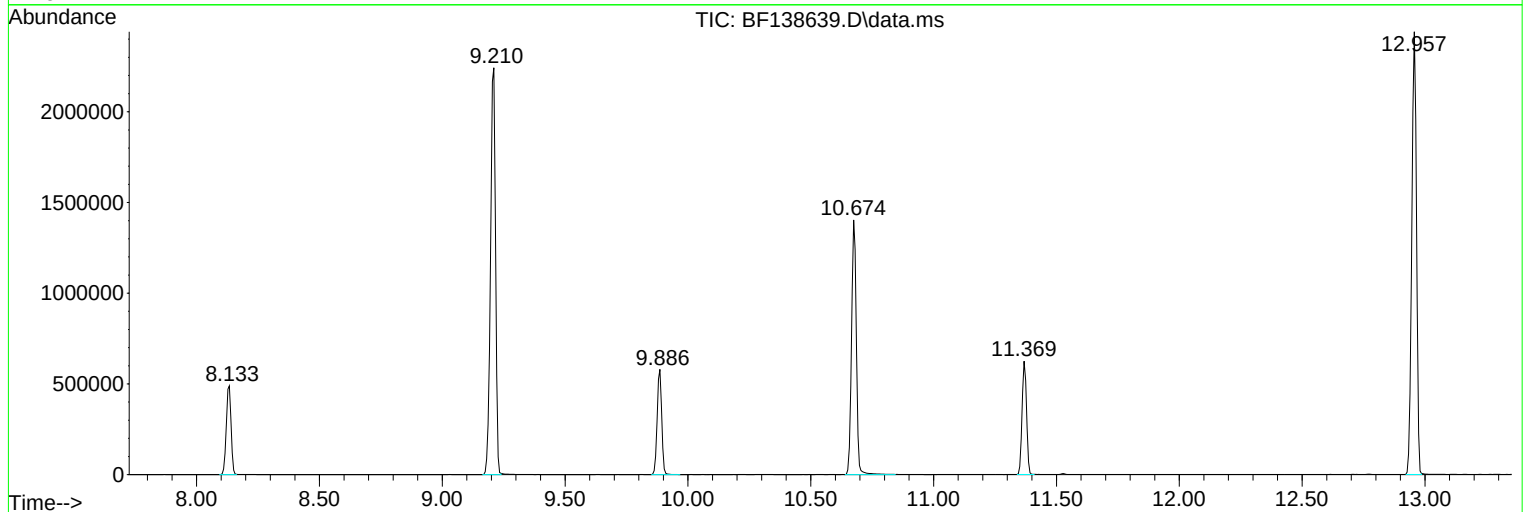
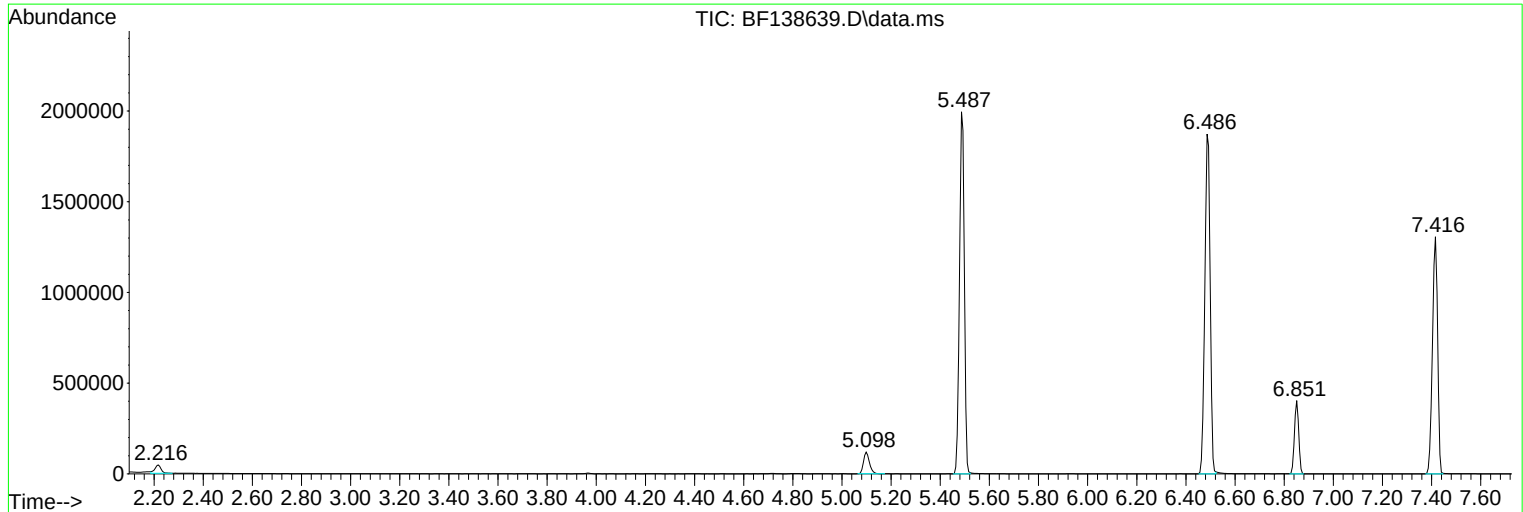
Sum of corrected areas: 19638271

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.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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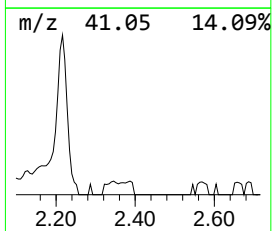
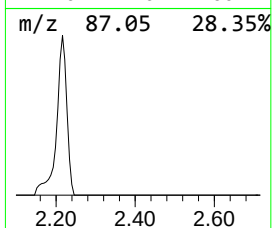
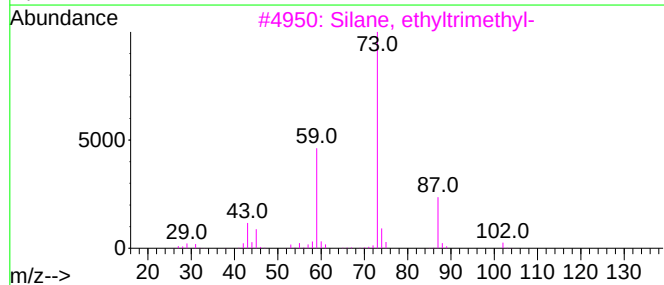
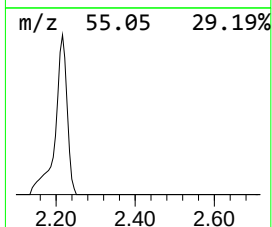
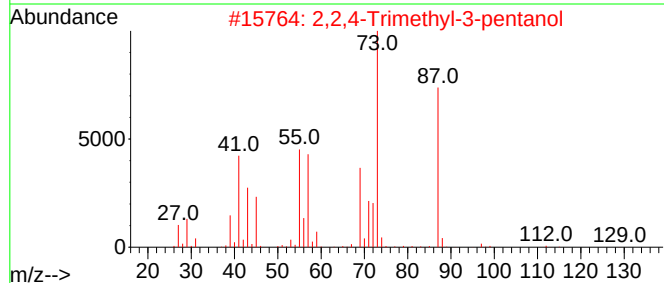
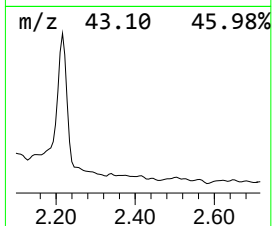
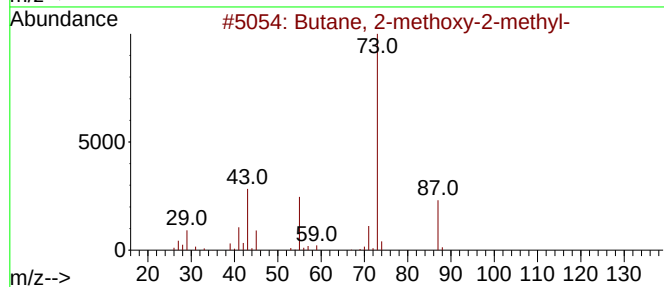
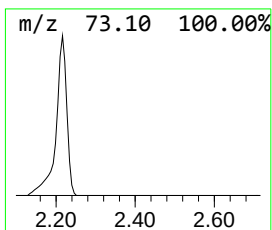
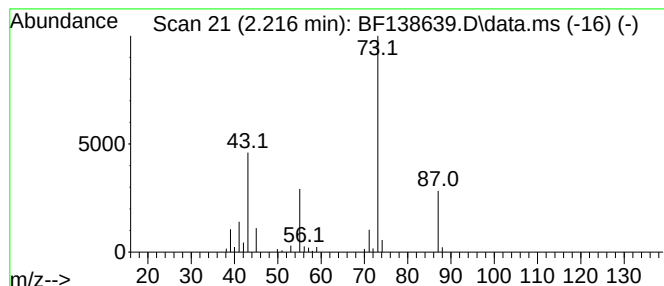
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.216	3.21 ng	77946	1,4-Dichlorobenzene-d4	6.851

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	39
3			Silane, ethyltrimethyl-	102	C5H14Si	003439-38-1	33
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	17
5			N-Ethylformamide	73	C3H7NO	000627-45-2	9



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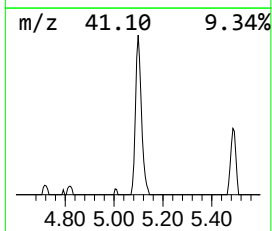
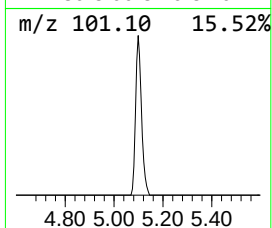
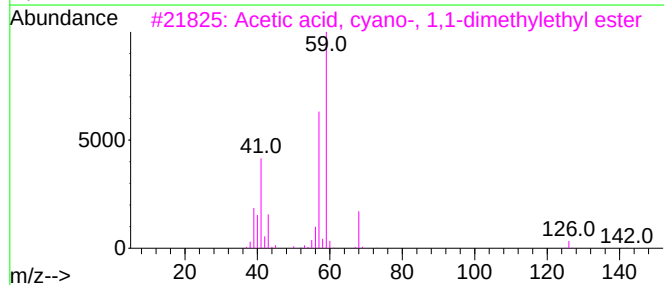
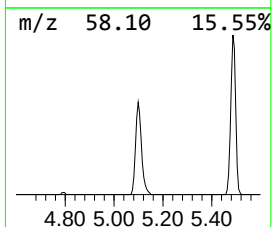
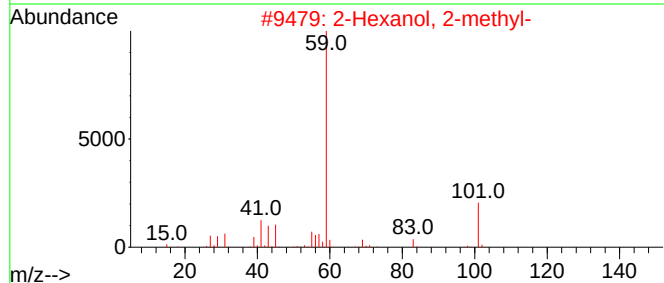
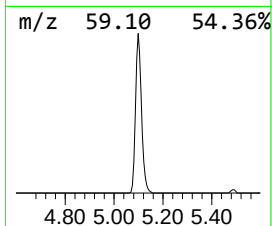
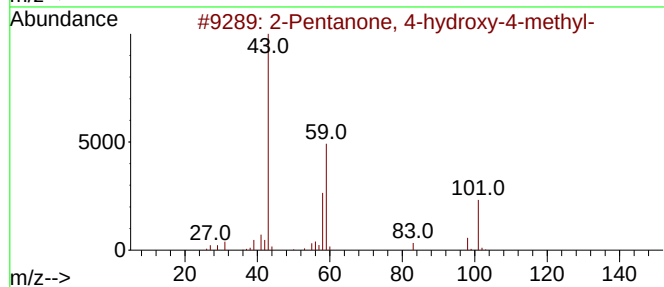
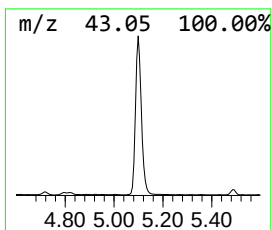
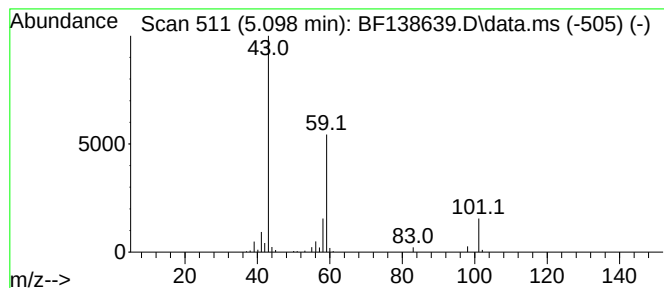
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.098	8.26 ng	200773	1,4-Dichlorobenzene-d4			6.851
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-		116	C6H12O2	000123-42-2	50
2	2-Hexanol, 2-methyl-		116	C7H16O	000625-23-0	33
3	Acetic acid, cyano-, 1,1-dimethy...		141	C7H11NO2	001116-98-9	25
4	1-Propen-2-ol, acetate		100	C5H8O2	000108-22-5	12
5	3-Hexanol, 4-methyl-		116	C7H16O	000615-29-2	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	2.216	3.2	ng	77946	1	6.851	486354	20.0
2-Pentanone, 4-...	5.098	8.3	ng	200773	1	6.851	486354	20.0