

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020122\
 Data File : BP008980.D
 Acq On : 01 Feb 2022 11:21
 Operator : CG/JU
 Sample : PB142379BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB142379BL

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_P\Methods\8270E-BP011422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008980.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.040	4	9	14	rVB	140967	177004	1.72%	0.349%
2	5.416	407	413	431	rBV	3554322	5556211	54.06%	10.970%
3	7.010	676	684	703	rBV	3091376	5353811	52.09%	10.571%
4	7.828	816	823	839	rVB	732998	1253191	12.19%	2.474%
5	8.987	1011	1020	1046	rBV	2060355	3641500	35.43%	7.190%
6	10.628	1290	1299	1318	rBV	849834	1617352	15.73%	3.193%
7	13.087	1710	1717	1743	rBV	4933419	8406112	81.78%	16.597%
8	14.469	1945	1952	1972	rBV	1109293	1914564	18.63%	3.780%
9	15.963	2199	2206	2236	rBV	3028316	5387799	52.42%	10.638%
10	17.222	2414	2420	2436	rBV	1241815	2018821	19.64%	3.986%
11	19.786	2850	2856	2869	rBV	8461190	10278707	100.00%	20.295%
12	21.316	3111	3116	3126	rBV	1733439	2365624	23.01%	4.671%
13	23.721	3519	3525	3539	rVB2	1209631	2676964	26.04%	5.285%

Sum of corrected areas: 50647660

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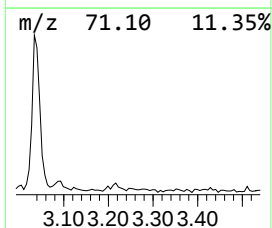
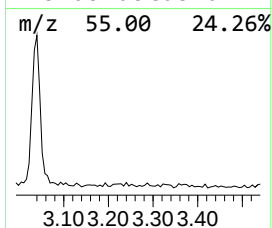
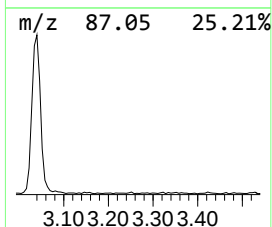
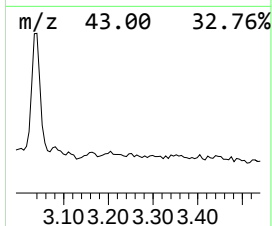
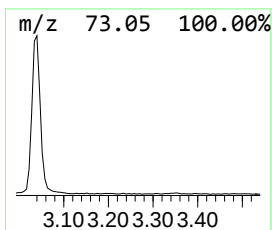
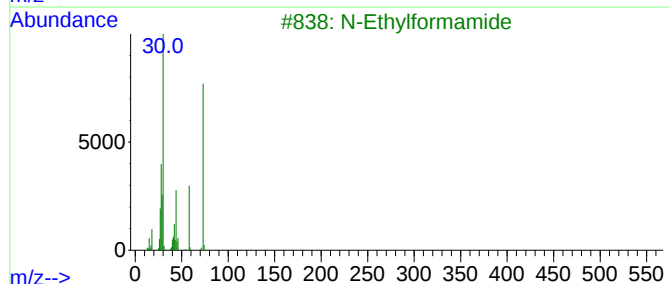
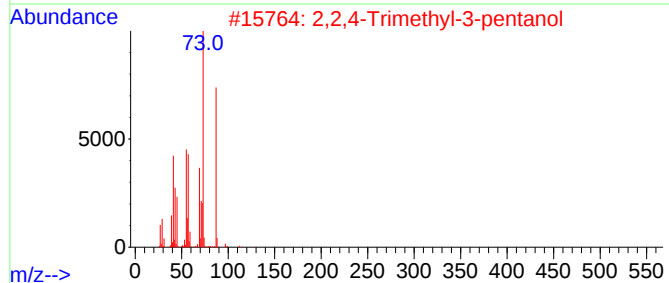
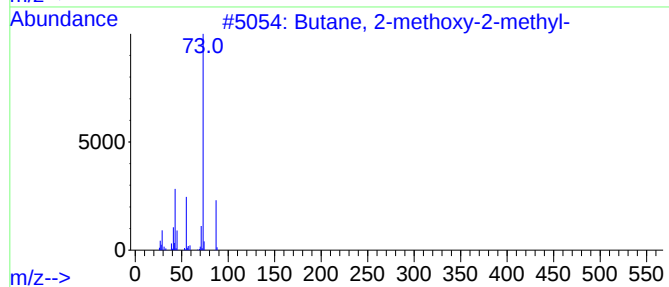
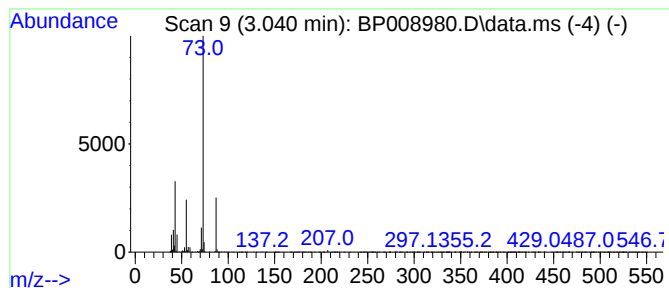
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	2.82 ng	177004	1,4-Dichlorobenzene-d4	7.828

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			N-Ethylformamide	73	C3H7NO	000627-45-2	9
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	9
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.040	2.8	ng	177004	1	7.828	1253190	20.0