

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
 Data File : BP008089.D
 Acq On : 23 Nov 2021 10:14
 Operator : CG/JU
 Sample : PB140937BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB140937BL

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-BP111921.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008089.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	14	rVB	119180	137996	1.23%	0.277%
2	5.411	405	412	427	rBV	3277315	4842362	43.27%	9.713%
3	7.005	675	683	699	rBV	3054280	5057188	45.19%	10.144%
4	7.822	814	822	830	rBV	609154	954415	8.53%	1.914%
5	8.981	1010	1019	1038	rBV	1925910	3237432	28.93%	6.494%
6	10.616	1289	1297	1307	rBV	798778	1341461	11.99%	2.691%
7	13.087	1707	1717	1734	rBV	4561061	7001150	62.57%	14.043%
8	14.463	1944	1951	1962	rBV	1284828	1925884	17.21%	3.863%
9	15.957	2197	2205	2226	rBV	3635923	5768870	51.55%	11.571%
10	17.222	2413	2420	2433	rBV	1691896	2551707	22.80%	5.118%
11	19.792	2851	2857	2869	rBV	8625835	11190073	100.00%	22.445%
12	21.316	3111	3116	3124	rBV	2171956	2940547	26.28%	5.898%
13	23.716	3517	3524	3533	rVB2	1421137	2907138	25.98%	5.831%

Sum of corrected areas: 49856223

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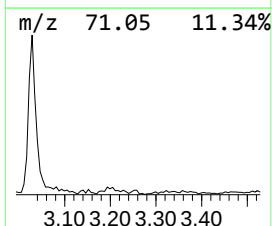
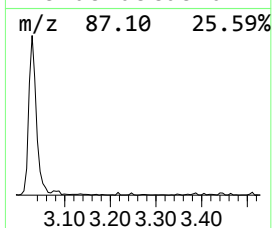
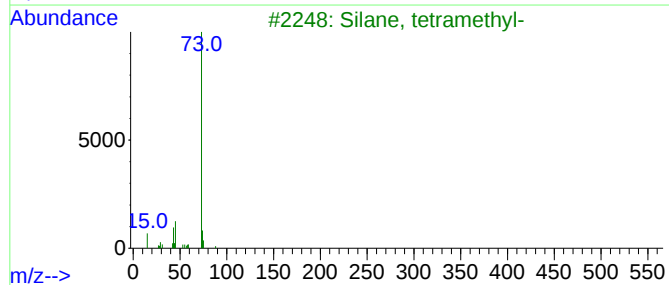
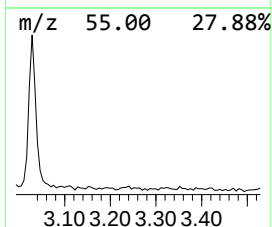
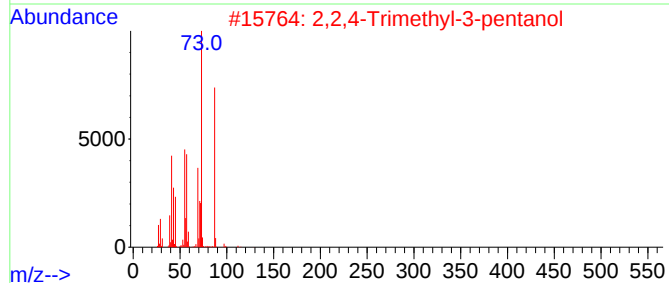
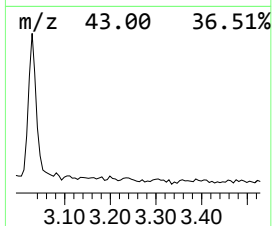
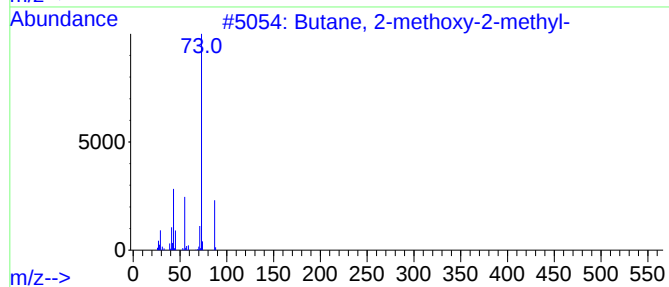
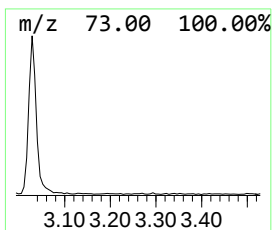
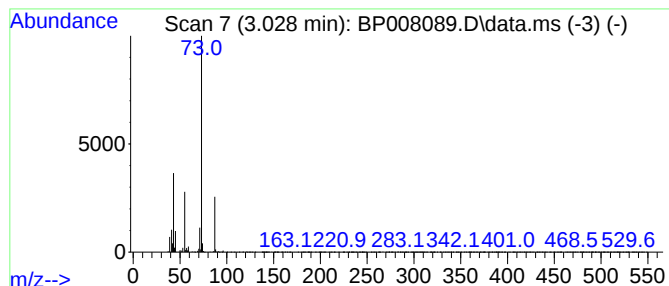
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
3.028	2.89 ng	137996	1,4-Dichlorobenzene-d4	7.822

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			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	12



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-metho...	3.028	2.9	ng	137996	1	7.822	954415	20.0