

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030376.D
 Acq On : 10 Jun 2021 16:10
 Operator : CG/JU
 Sample : M2613-17
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(98-102)

Manual Integrations
 APPROVED

mohammad
 6/11/2021 2:11:55 PM

Quant Time: Jun 10 17:01:28 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 08 20:51:08 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.845	152	278568	20.000	ng	0.00
21) Naphthalene-d8	10.633	136	1127348	20.000	ng	# 0.00
39) Acenaphthene-d10	14.468	164	791332	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1894133	20.000	ng	0.00
76) Chrysene-d12	21.362	240	1915409	20.000	ng	0.00
86) Perylene-d12	23.644	264	1662101	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	2242642	130.875	ng	0.00
7) Phenol-d6	7.010	99	2697967	108.762	ng	0.00
23) Nitrobenzene-d5	9.004	82	1849679	79.933	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	1555840	156.378	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	4654514	77.350	ng	0.00
79) Terphenyl-d14	19.821	244	9088071	83.338	ng	0.00
Target Compounds						
32) Benzoic acid	9.898	122	5956m	5.389	ng	Qvalue
84) Bis(2-ethylhexyl)phtha...	21.280	149	21224	4.488	ng	92

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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