

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061421\
 Data File : BM030438.D
 Acq On : 14 Jun 2021 15:56
 Operator : CG/JU
 Sample : M2625-30RE
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(174-178)RE

Manual Integrations
 APPROVED

mohammad
 6/15/2021 2:40:29 PM

Quant Time: Jun 14 18:01:09 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 14:10:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	178732	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	644970	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	379053	20.000	ng	0.00
64) Phenanthrene-d10	17.192	188	768486	20.000	ng	0.00
76) Chrysene-d12	21.356	240	861411	20.000	ng	0.00
86) Perylene-d12	23.633	264	905515	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1414801	128.683	ng	0.02
7) Phenol-d6	7.016	99	1477775	92.849	ng	0.02
23) Nitrobenzene-d5	8.987	82	1202534	90.833	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	752649	157.929	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	2701562	93.726	ng	0.00
79) Terphenyl-d14	19.809	244	4339017	88.474	ng	0.00
Target Compounds						Qvalue
32) Benzoic acid	9.940	122	3685m	5.421	ng	

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

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