

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061021\
 Data File : BM030380.D
 Acq On : 10 Jun 2021 18:38
 Operator : CG/JU
 Sample : M2613-21
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 18-B12(114-118)

Manual Integrations
 APPROVED

mohammad
 6/11/2021 2:11:59 PM

Quant Time: Jun 11 07:53:18 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	7.857	152	1019	20.000	ng	0.00
21) Naphthalene-d8	10.645	136	2785	20.000	ng	# 0.00
39) Acenaphthene-d10	14.474	164	1144	20.000	ng	# 0.00
64) Phenanthrene-d10	17.210	188	1574m	20.000	ng	0.00
76) Chrysene-d12	21.374	240	2770m	20.000	ng	0.00
86) Perylene-d12	23.656	264	2281	20.000	ng	# 0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	163	2.600	ng	0.00
7) Phenol-d6	7.028	99	10445	115.108	ng	0.02
23) Nitrobenzene-d5	9.004	82	1764	30.857	ng	0.00
42) 2,4,6-Tribromophenol	0.000	330	0	0.000	ng	
45) 2-Fluorobiphenyl	13.098	172	6912	79.455	ng	0.00
79) Terphenyl-d14	19.827	244	9591	60.816	ng	0.00
Target Compounds						
81) Benzo(a)anthracene	21.368	228	1344	6.667	ng	# 72
83) Chrysene	21.415	228	2784	14.210	ng	# 90

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

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