

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060921\
 Data File : BM030347.D
 Acq On : 09 Jun 2021 18:03
 Operator : CG/JU
 Sample : PB136946BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB136946BL

Quant Time: Jun 09 18:52:59 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	7.851	152	189390	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	745932	20.000	ng	# 0.00
39) Acenaphthene-d10	14.474	164	491351	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1120152	20.000	ng	# 0.00
76) Chrysene-d12	21.368	240	1243661	20.000	ng	0.00
86) Perylene-d12	23.650	264	1311474	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1781878	152.950	ng	0.00
7) Phenol-d6	7.010	99	2289101	135.731	ng	0.00
23) Nitrobenzene-d5	9.004	82	1350367	88.194	ng	0.00
42) 2,4,6-Tribromophenol	15.957	330	917406	148.504	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	3292239	88.114	ng	0.00
79) Terphenyl-d14	19.821	244	6338194	89.515	ng	0.00

Target Compounds	Qvalue

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

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