

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061522\
 Data File : BM035476.D
 Acq On : 15 Jun 2022 16:58
 Operator : CG/JU
 Sample : PB145557BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB145557BL

Quant Time: Jun 15 23:48:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM061022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 10 16:05:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	250369	20.000	ng	0.00
21) Naphthalene-d8	10.616	136	1077148	20.000	ng	0.00
39) Acenaphthene-d10	14.457	164	688053	20.000	ng	0.00
64) Phenanthrene-d10	17.203	188	1493244	20.000	ng	# 0.00
76) Chrysene-d12	21.391	240	1441752	20.000	ng	0.00
86) Perylene-d12	23.738	264	1447865	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.404	112	2182328	147.812	ng	0.00
7) Phenol-d6	7.004	99	2784646	136.104	ng	0.00
23) Nitrobenzene-d5	8.980	82	1645303	83.007	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1053422	143.803	ng	0.00
45) 2-Fluorobiphenyl	13.080	172	4068802	84.481	ng	0.00
79) Terphenyl-d14	19.833	244	6256041	82.850	ng	0.00

Target Compounds	Qvalue

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

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