

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072022\
 Data File : BM035880.D
 Acq On : 20 Jul 2022 14:27
 Operator : CG/JU
 Sample : PB145957BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB145957BL

Quant Time: Jul 20 15:42:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 18 17:08:15 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	349191	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1407084	20.000	ng	0.00
39) Acenaphthene-d10	14.533	164	764299	20.000	ng	0.00
64) Phenanthrene-d10	17.269	188	1347870	20.000	ng	# 0.00
76) Chrysene-d12	21.445	240	895466	20.000	ng	0.00
86) Perylene-d12	23.821	264	828139	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.458	112	3110896	138.114	ng	0.00
7) Phenol-d6	7.063	99	3763459	132.873	ng	0.00
23) Nitrobenzene-d5	9.063	82	2247639	85.433	ng	0.00
42) 2,4,6-Tribromophenol	16.016	330	982976	114.980	ng	0.00
45) 2-Fluorobiphenyl	13.163	172	4792044	88.625	ng	0.00
79) Terphenyl-d14	19.898	244	4908378	106.649	ng	0.00

Target Compounds	Qvalue

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

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