

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071922\
 Data File : BM035871.D
 Acq On : 19 Jul 2022 19:19
 Operator : CG/JU
 Sample : PB145925BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB145925BL

Quant Time: Jul 20 02:46:22 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	7.898	152	323535	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	1284677	20.000	ng	0.00
39) Acenaphthene-d10	14.533	164	665107	20.000	ng	0.00
64) Phenanthrene-d10	17.268	188	1177048	20.000	ng	0.00
76) Chrysene-d12	21.445	240	858069	20.000	ng	0.00
86) Perylene-d12	23.821	264	845919	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.463	112	2801796	134.255	ng	0.00
7) Phenol-d6	7.069	99	3337738	127.187	ng	0.00
23) Nitrobenzene-d5	9.069	82	2011625	83.748	ng	0.00
42) 2,4,6-Tribromophenol	16.015	330	837425	112.563	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	4162164	88.456	ng	0.00
79) Terphenyl-d14	19.898	244	4466523	101.278	ng	0.00

Target Compounds	Qvalue

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

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