

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052022\
 Data File : BM035171.D
 Acq On : 20 May 2022 11:46
 Operator : CG/JU
 Sample : PB144941BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB144941BL

Quant Time: May 23 04:16:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042722.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 20 04:53:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	158405	20.000	ng	0.00
21) Naphthalene-d8	10.704	136	652694	20.000	ng	0.00
39) Acenaphthene-d10	14.545	164	426039	20.000	ng	0.00
64) Phenanthrene-d10	17.286	188	858382	20.000	ng	0.00
76) Chrysene-d12	21.468	240	636070	20.000	ng	# 0.00
86) Perylene-d12	23.856	264	577857	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.481	112	894168	99.587	ng	0.00
7) Phenol-d6	7.075	99	1105067	95.338	ng	0.00
23) Nitrobenzene-d5	9.063	82	1062635	87.919	ng	0.00
42) 2,4,6-Tribromophenol	16.033	330	457837	98.790	ng	0.00
45) 2-Fluorobiphenyl	13.169	172	2627742	89.537	ng	0.00
79) Terphenyl-d14	19.915	244	3447510	104.943	ng	0.00

Target Compounds	Qvalue

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

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