

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080822\
 Data File : BM036150.D
 Acq On : 08 Aug 2022 18:40
 Operator : CG/JU
 Sample : PB146776BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146776BL

Quant Time: Aug 08 23:39:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 08 23:36:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	331456	20.000	ng	0.00
21) Naphthalene-d8	10.651	136	1446809	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	923481	20.000	ng	0.00
64) Phenanthrene-d10	17.233	188	1934730	20.000	ng	0.00
76) Chrysene-d12	21.415	240	1667804	20.000	ng	0.00
86) Perylene-d12	23.768	264	1539573	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	2817596	137.818	ng	0.00
7) Phenol-d6	7.058	99	3602240	138.473	ng	0.00
23) Nitrobenzene-d5	9.016	82	2222440	85.986	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	1461355	134.890	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	5208592	83.276	ng	0.00
79) Terphenyl-d14	19.863	244	7989287	94.486	ng	0.00

Target Compounds	Qvalue

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

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