

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080922\
 Data File : BM036169.D
 Acq On : 09 Aug 2022 11:06
 Operator : CG/JU
 Sample : PB146619BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146619BL

Quant Time: Aug 10 03:30:29 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	7.839	152	297654	20.000	ng	0.00
21) Naphthalene-d8	10.645	136	1303923	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	859511	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	1910743	20.000	ng	0.00
76) Chrysene-d12	21.409	240	1860828	20.000	ng	-0.01
86) Perylene-d12	23.762	264	1917516	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	2615959	142.486	ng	0.00
7) Phenol-d6	7.057	99	3349276	143.370	ng	0.00
23) Nitrobenzene-d5	9.016	82	2050467	88.026	ng	0.00
42) 2,4,6-Tribromophenol	15.980	330	1503817	149.141	ng	0.00
45) 2-Fluorobiphenyl	13.116	172	4918249	84.486	ng	0.00
79) Terphenyl-d14	19.856	244	8610364	91.268	ng	0.00

Target Compounds	Qvalue

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

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