

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080422\
 Data File : BM036094.D
 Acq On : 04 Aug 2022 12:11
 Operator : CG/JU
 Sample : PB146679BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146679BL

Quant Time: Aug 05 03:38:37 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 03:36:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.857	152	273825	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	1154007	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	679983	20.000	ng	0.00
64) Phenanthrene-d10	17.239	188	1310115	20.000	ng	0.00
76) Chrysene-d12	21.421	240	1037691	20.000	ng	0.00
86) Perylene-d12	23.786	264	982098	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	2305150	136.483	ng	0.00
7) Phenol-d6	7.034	99	2848956	132.566	ng	0.00
23) Nitrobenzene-d5	9.028	82	1721170	83.488	ng	0.00
42) 2,4,6-Tribromophenol	15.986	330	1040300	130.411	ng	0.00
45) 2-Fluorobiphenyl	13.128	172	3915453	85.018	ng	0.00
79) Terphenyl-d14	19.868	244	5159325	98.068	ng	0.00

Target Compounds	Qvalue

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

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