

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081022\
 Data File : BM036188.D
 Acq On : 10 Aug 2022 12:31
 Operator : CG/JU
 Sample : PB146851BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146851BL

Quant Time: Aug 10 22:22:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 10 22:19:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	240263	20.000	ng	0.00
21) Naphthalene-d8	10.639	136	961019	20.000	ng	0.00
39) Acenaphthene-d10	14.474	164	519572	20.000	ng	0.00
64) Phenanthrene-d10	17.215	188	968663	20.000	ng	0.00
76) Chrysene-d12	21.398	240	800771	20.000	ng	0.00
86) Perylene-d12	23.744	264	806139	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.422	112	2098078	141.575	ng	0.00
7) Phenol-d6	7.040	99	2567583	136.162	ng	0.00
23) Nitrobenzene-d5	9.004	82	1565060	91.161	ng	0.00
42) 2,4,6-Tribromophenol	15.968	330	756900	124.178	ng	0.00
45) 2-Fluorobiphenyl	13.104	172	3246430	92.254	ng	0.00
79) Terphenyl-d14	19.845	244	3929128	96.781	ng	0.00

Target Compounds	Qvalue

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

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