

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080122\
 Data File : BM036037.D
 Acq On : 01 Aug 2022 19:18
 Operator : CG/JU
 Sample : PB146582BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB146582BL

Quant Time: Aug 02 07:44:50 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 01 17:50:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	323769	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1341118	20.000	ng	0.00
39) Acenaphthene-d10	14.510	164	758844	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1439491	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1091412	20.000	ng	0.00
86) Perylene-d12	23.791	264	989074	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	2875396	143.984	ng	0.00
7) Phenol-d6	7.045	99	3503868	137.890	ng	0.00
23) Nitrobenzene-d5	9.039	82	2204091	91.997	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1150353	129.221	ng	0.00
45) 2-Fluorobiphenyl	13.139	172	4822119	93.823	ng	0.00
79) Terphenyl-d14	19.874	244	6019012	108.778	ng	0.00

Target Compounds	Qvalue

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

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