

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092023\
 Data File : BM041932.D
 Acq On : 20 Sep 2023 10:35
 Operator : MA/JU
 Sample : PB155666BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB155666BL

Quant Time: Sep 20 12:11:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM091823.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 19 04:49:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	299511	20.000	ng	-0.01
21) Naphthalene-d8	10.775	136	1218665	20.000	ng	-0.01
39) Acenaphthene-d10	14.604	164	774735	20.000	ng	-0.01
64) Phenanthrene-d10	17.357	188	1664029	20.000	ng	-0.01
76) Chrysene-d12	21.527	240	1451698	20.000	ng	-0.02
86) Perylene-d12	23.944	264	1460078	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	2644022	147.036	ng	0.00
7) Phenol-d6	7.110	99	3482064	139.463	ng	0.00
23) Nitrobenzene-d5	9.151	82	2139148	89.373	ng	-0.01
42) 2,4,6-Tribromophenol	16.098	330	1327999	137.384	ng	-0.01
45) 2-Fluorobiphenyl	13.227	172	4682936	88.720	ng	0.00
79) Terphenyl-d14	19.968	244	7381582	92.609	ng	0.00

Target Compounds	Qvalue

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

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