

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080922\
 Data File : BM036173.D
 Acq On : 09 Aug 2022 13:45
 Operator : CG/JU
 Sample : N3962-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-197

Quant Time: Aug 10 03:32:56 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	218041	20.000	ng	0.00
21) Naphthalene-d8	10.645	136	829090	20.000	ng	0.00
39) Acenaphthene-d10	14.486	164	463039	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	948703	20.000	ng	0.00
76) Chrysene-d12	21.409	240	1009852	20.000	ng	-0.01
86) Perylene-d12	23.768	264	1111363	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1226320	91.184	ng	0.00
7) Phenol-d6	7.051	99	1470433	85.926	ng	-0.01
23) Nitrobenzene-d5	9.010	82	895876	60.486	ng	-0.01
42) 2,4,6-Tribromophenol	15.974	330	531134	97.778	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	1898062	60.523	ng	-0.01
79) Terphenyl-d14	19.850	244	3045453	59.484	ng	-0.01
Target Compounds						
78) Pyrene	19.645	202	133968	2.139	ng	Qvalue 99

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

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