

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080822\
 Data File : BM036157.D
 Acq On : 08 Aug 2022 23:03
 Operator : CG/JU
 Sample : N3960-14
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RVE-161

Quant Time: Aug 08 23:44:01 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	7.846	152	289174	20.00	ng	0.00
21) Naphthalene-d8	10.645	136	1128345	20.00	ng	0.00
39) Acenaphthene-d10	14.486	164	620625	20.00	ng	0.00
64) Phenanthrene-d10	17.227	188	1209928	20.00	ng	0.00
76) Chrysene-d12	21.409	240	1148850	20.00	ng	-0.01
86) Perylene-d12	23.762	264	1239812	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1039404	58.27	ng	0.00
7) Phenol-d6	7.051	99	1311912	57.80	ng	-0.01
23) Nitrobenzene-d5	9.016	82	903269	44.81	ng	0.00
42) 2,4,6-Tribromophenol	15.974	330	467984	64.28	ng	-0.01
45) 2-Fluorobiphenyl	13.110	172	1949374	46.38	ng	-0.01
79) Terphenyl-d14	19.851	244	2655324	45.59	ng	-0.01
Target Compounds						
78) Pyrene	19.645	202	185081	2.60	ng	Qvalue 99

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

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