

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080222\
 Data File : BM036055.D
 Acq On : 02 Aug 2022 17:10
 Operator : CG/JU
 Sample : N3976-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP1-0729

Quant Time: Aug 03 00:53:03 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	391616	20.000	ng	0.00
21) Naphthalene-d8	10.675	136	1605118	20.000	ng	0.00
39) Acenaphthene-d10	14.504	164	941387	20.000	ng	0.00
64) Phenanthrene-d10	17.245	188	1826318	20.000	ng	0.00
76) Chrysene-d12	21.427	240	1370427	20.000	ng	0.00
86) Perylene-d12	23.791	264	1159925	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	3495090	144.694	ng	0.01
7) Phenol-d6	7.045	99	3959037	128.810	ng	0.00
23) Nitrobenzene-d5	9.039	82	2864266	99.888	ng	0.00
42) 2,4,6-Tribromophenol	15.992	330	1699937	153.928	ng	0.00
45) 2-Fluorobiphenyl	13.139	172	6236145	97.808	ng	0.00
79) Terphenyl-d14	19.874	244	8328578	119.872	ng	0.00

Target Compounds	Qvalue

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

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