

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081122\
 Data File : BM036219.D
 Acq On : 11 Aug 2022 21:12
 Operator : CG/JU
 Sample : N4082-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 CRUSHED-MASONARY-PILE-A

Quant Time: Aug 12 02:19:24 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 10 22:19:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	262490	20.000	ng	-0.01
21) Naphthalene-d8	10.628	136	1145693	20.000	ng	-0.01
39) Acenaphthene-d10	14.468	164	722712	20.000	ng	0.00
64) Phenanthrene-d10	17.209	188	1538723	20.000	ng	0.00
76) Chrysene-d12	21.392	240	1587959	20.000	ng	0.00
86) Perylene-d12	23.733	264	1535049	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.410	112	361524	22.329	ng	0.00
7) Phenol-d6	7.022	99	1266583	61.481	ng	-0.01
23) Nitrobenzene-d5	8.998	82	1339421	65.442	ng	0.00
42) 2,4,6-Tribromophenol	15.963	330	27765	3.275	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	3228366	65.954	ng	0.00
79) Terphenyl-d14	19.839	244	5603678	69.605	ng	0.00

Target Compounds	Qvalue

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

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