

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082123\
 Data File : BM041444.D
 Acq On : 21 Aug 2023 12:35
 Operator : MA/JU
 Sample : 04080-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETGI-340

Quant Time: Aug 22 03:35:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 22 03:32:15 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.757	152	131205	20.000	ng	0.00
21) Naphthalene-d8	10.557	136	521672	20.000	ng	0.00
39) Acenaphthene-d10	14.415	164	358910	20.000	ng	0.00
64) Phenanthrene-d10	17.180	188	851388	20.000	ng	0.00
76) Chrysene-d12	21.386	240	935706	20.000	ng	0.00
86) Perylene-d12	23.727	264	1135848	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.345	112	878968	111.446	ng	0.00
7) Phenol-d6	6.957	99	1029348	96.115	ng	0.00
23) Nitrobenzene-d5	8.933	82	758991	69.452	ng	0.00
42) 2,4,6-Tribromophenol	15.921	330	616282	108.201	ng	0.00
45) 2-Fluorobiphenyl	13.033	172	1839799	68.690	ng	0.00
79) Terphenyl-d14	19.815	244	3702258	68.852	ng	0.00

Target Compounds	Qvalue

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

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