

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081621\
 Data File : BM031773.D
 Acq On : 17 Aug 2021 04:54
 Operator : CG/JU
 Sample : M3407-27
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 29N

Quant Time: Aug 17 07:28:57 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 16:42:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.551	152	48162	20.000	ng	0.00
21) Naphthalene-d8	10.310	136	197723	20.000	ng	0.00
39) Acenaphthene-d10	14.186	164	125882	20.000	ng	0.00
64) Phenanthrene-d10	16.939	188	268681	20.000	ng	0.00
76) Chrysene-d12	21.133	240	257004	20.000	ng	-0.01
86) Perylene-d12	23.286	264	219809	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	442634	147.508	ng	0.00
7) Phenol-d6	6.746	99	514943	118.428	ng	0.00
23) Nitrobenzene-d5	8.704	82	448403	94.957	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	224808	145.000	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	890886	94.198	ng	0.00
79) Terphenyl-d14	19.586	244	1552899	101.805	ng	0.00

Target Compounds	Qvalue

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

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