

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081621\
 Data File : BM031772.D
 Acq On : 17 Aug 2021 04:17
 Operator : CG/JU
 Sample : M3407-06
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 25S

Quant Time: Aug 17 07:28:44 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	50839	20.000	ng	0.00
21) Naphthalene-d8	10.310	136	221871	20.000	ng	# 0.00
39) Acenaphthene-d10	14.186	164	147082	20.000	ng	0.00
64) Phenanthrene-d10	16.939	188	309506	20.000	ng	0.00
76) Chrysene-d12	21.139	240	271626	20.000	ng	0.00
86) Perylene-d12	23.286	264	224632	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	439099	138.624	ng	0.00
7) Phenol-d6	6.745	99	512269	111.609	ng	0.00
23) Nitrobenzene-d5	8.704	82	458437	86.516	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	241177	133.136	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	953914	86.324	ng	0.00
79) Terphenyl-d14	19.586	244	1695271	105.156	ng	0.00

Target Compounds	Qvalue

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

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