

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081621\
 Data File : BM031761.D
 Acq On : 16 Aug 2021 21:29
 Operator : CG/JU
 Sample : M3407-18
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 27S

Quant Time: Aug 17 03:10:13 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	36164	20.000	ng	0.00
21) Naphthalene-d8	10.310	136	144683	20.000	ng	# 0.00
39) Acenaphthene-d10	14.186	164	94744	20.000	ng	0.00
64) Phenanthrene-d10	16.939	188	208534	20.000	ng	0.00
76) Chrysene-d12	21.139	240	212601	20.000	ng	0.00
86) Perylene-d12	23.292	264	204760	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.193	112	336788	149.470	ng	0.00
7) Phenol-d6	6.751	99	369080	113.043	ng	0.00
23) Nitrobenzene-d5	8.704	82	351895	101.838	ng	0.00
42) 2,4,6-Tribromophenol	15.692	330	180726	154.877	ng	0.00
45) 2-Fluorobiphenyl	12.804	172	697274	97.957	ng	0.00
79) Terphenyl-d14	19.592	244	1380371	109.395	ng	0.00

Target Compounds	Qvalue

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

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