

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
 Data File : BM031808.D
 Acq On : 18 Aug 2021 16:32
 Operator : CG/JU
 Sample : M3429-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-1

Quant Time: Aug 18 18:10:21 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 18 14:13:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.540	152	29800	20.000	ng	# 0.00
21) Naphthalene-d8	10.298	136	110452	20.000	ng	# 0.00
39) Acenaphthene-d10	14.174	164	68345	20.000	ng	0.00
64) Phenanthrene-d10	16.933	188	144264	20.000	ng	0.00
76) Chrysene-d12	21.133	240	142330	20.000	ng	0.00
86) Perylene-d12	23.280	264	142623	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.181	112	281481	151.603	ng	0.00
7) Phenol-d6	6.734	99	312631	116.202	ng	0.00
23) Nitrobenzene-d5	8.692	82	258955	98.167	ng	0.00
42) 2,4,6-Tribromophenol	15.680	330	115928	137.721	ng	0.00
45) 2-Fluorobiphenyl	12.792	172	492137	95.843	ng	0.00
79) Terphenyl-d14	19.580	244	870343	103.029	ng	0.00

Target Compounds	Qvalue

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

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