

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
 Data File : BM031813.D
 Acq On : 18 Aug 2021 19:39
 Operator : CG/JU
 Sample : M3449-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP3

Quant Time: Aug 19 03:05:03 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	23532	20.000	ng	# 0.00
21) Naphthalene-d8	10.298	136	88539	20.000	ng	# 0.00
39) Acenaphthene-d10	14.175	164	55106	20.000	ng	0.00
64) Phenanthrene-d10	16.933	188	118379	20.000	ng	0.00
76) Chrysene-d12	21.133	240	119060	20.000	ng	0.00
86) Perylene-d12	23.274	264	131939	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.175	112	203935	139.094	ng	0.00
7) Phenol-d6	6.734	99	216080	101.708	ng	0.00
23) Nitrobenzene-d5	8.693	82	202404	95.719	ng	0.00
42) 2,4,6-Tribromophenol	15.680	330	89045	131.199	ng	0.00
45) 2-Fluorobiphenyl	12.792	172	380519	91.909	ng	0.00
79) Terphenyl-d14	19.580	244	654119	92.567	ng	0.00

Target Compounds	Qvalue

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

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