

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081821\
 Data File : BM031806.D
 Acq On : 18 Aug 2021 15:18
 Operator : CG/JU
 Sample : M3424-12
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 19S

Quant Time: Aug 18 15:59:04 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	24065	20.000	ng	# 0.00
21) Naphthalene-d8	10.298	136	95359	20.000	ng	# 0.00
39) Acenaphthene-d10	14.174	164	61246	20.000	ng	0.00
64) Phenanthrene-d10	16.933	188	131451	20.000	ng	0.00
76) Chrysene-d12	21.133	240	138226	20.000	ng	0.00
86) Perylene-d12	23.280	264	151564	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.175	112	31510	21.015	ng	0.00
7) Phenol-d6	6.734	99	51855	23.867	ng	0.00
23) Nitrobenzene-d5	8.692	82	139202	61.122	ng	0.00
42) 2,4,6-Tribromophenol	15.680	330	41095	54.479	ng	0.00
45) 2-Fluorobiphenyl	12.792	172	275327	59.835	ng	0.00
79) Terphenyl-d14	19.580	244	510376	62.211	ng	0.00

Target Compounds Qvalue

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

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