

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081719\
 Data File : BG042270.D
 Acq On : 16 Aug 2019 21:47
 Operator : HP/JU
 Sample : K4231-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 REACTOR-3-B-(0-3)

Quant Time: Aug 17 00:36:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG072719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 08:57:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	58122	20.00	ng	-0.03
21) Naphthalene-d8	10.84	136	218396	20.00	ng	-0.03
39) Acenaphthene-d10	14.67	164	156721	20.00	ng	-0.03
64) Phenanthrene-d10	17.42	188	382638	20.00	ng	-0.03
76) Chrysene-d12	21.71	240	449689	20.00	ng	-0.03
87) Perylene-d12	24.96	264	521945	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	386472	129.37	ng	-0.02
7) Phenol-d6	7.18	99	464827	115.42	ng	-0.02
23) Nitrobenzene-d5	9.18	82	319275	100.60	ng	-0.03
42) 2,4,6-Tribromophenol	16.17	330	343409	192.91	ng	-0.02
45) 2-Fluorobiphenyl	13.29	172	957619	107.77	ng	-0.03
79) Terphenyl-d14	20.04	244	1890863	96.22	ng	-0.03

Target Compounds Qvalue

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

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