

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081719\
 Data File : BF116351.D
 Acq On : 17 Aug 2019 12:30
 Operator : HP/JU
 Sample : K4377-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RBR-29

Quant Time: Aug 19 06:41:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	149722	20.00	ng	-0.04
21) Naphthalene-d8	8.09	136	571106	20.00	ng	-0.04
39) Acenaphthene-d10	9.84	164	299967	20.00	ng	-0.05
64) Phenanthrene-d10	11.33	188	513793	20.00	ng	-0.04
76) Chrysene-d12	13.97	240	362717	20.00	ng	-0.04
87) Perylene-d12	15.42	264	342934	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1100423	123.04	ng	-0.02
7) Phenol-d6	6.46	99	1405888	124.59	ng	-0.03
23) Nitrobenzene-d5	7.37	82	914934	92.47	ng	-0.04
42) 2,4,6-Tribromophenol	10.63	330	354416	121.87	ng	-0.04
45) 2-Fluorobiphenyl	9.16	172	1408524	87.95	ng	-0.04
79) Terphenyl-d14	12.91	244	1437375	83.50	ng	-0.04
Target Compounds						
10) Phenol	6.46	94	49404	3.718	ng	84
50) Dimethylphthalate	9.56	163	319114	15.624	ng	99

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

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