

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081619\
 Data File : BF116343.D
 Acq On : 17 Aug 2019 3:02
 Operator : HP/JU
 Sample : K4391-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BOTTOM-1

Quant Time: Aug 19 08:37:41 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	6.82	152	140113	20.00	ng	-0.03
21) Naphthalene-d8	8.10	136	535972	20.00	ng	-0.03
39) Acenaphthene-d10	9.85	164	274335	20.00	ng	-0.04
64) Phenanthrene-d10	11.33	188	432587	20.00	ng	-0.04
76) Chrysene-d12	13.97	240	281602	20.00	ng	-0.04
87) Perylene-d12	15.43	264	316485	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	767678	91.72	ng	-0.02
7) Phenol-d6	6.46	99	994745	94.20	ng	-0.02
23) Nitrobenzene-d5	7.38	82	600664	64.69	ng	-0.04
42) 2,4,6-Tribromophenol	10.64	330	220419	82.87	ng	-0.04
45) 2-Fluorobiphenyl	9.17	172	971106	66.30	ng	-0.04
79) Terphenyl-d14	12.92	244	821313	61.46	ng	-0.04
Target Compounds						Qvalue
10) Phenol	6.47	94	53997	4.342	ng	94
37) 2-Methylnaphthalene	8.81	142	134117	8.074	ng	98
38) 1-Methylnaphthalene	8.91	142	77554	4.935	ng	99
50) Dimethylphthalate	9.56	163	101713	5.445	ng	99
71) Phenanthrene	11.36	178	92407	4.179	ng	98
75) Fluoranthene	12.55	202	103746	4.481	ng	99
78) Pyrene	12.78	202	94445	4.449	ng	99
81) Benzo(a)anthracene	13.97	228	36817	2.046	ng	98
88) Benzo(b)fluoranthene	15.02	252	38236m	2.089	ng	

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

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