

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050818\
 Data File : BF105164.D
 Acq On : 9 May 2018 00:18
 Operator : JU/SJ
 Sample : J2133-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-WATER-01-QT2-2018

Quant Time: May 09 09:28:56 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 08 16:33:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	201146	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	804193	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	403137	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	788929	20.00	ng	0.00
75) Chrysene-d12	14.00	240	659538	20.00	ng	-0.02
86) Perylene-d12	15.54	264	595477	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1524462	124.86	ng	0.00
7) Phenol-d6	6.42	99	1824000	125.46	ng	0.00
23) Nitrobenzene-d5	7.36	82	1144106	96.42	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	597065	134.24	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	2069932	79.69	ng	0.00
78) Terphenyl-d14	12.91	244	2230021	78.54	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	8.93	237	55187	7.956	ng	98
53) 2,4-Dinitrophenol	9.87	184	11872	16.500	ng	# 24
69) Pentachlorophenol	11.11	266	45770	10.900	ng	94

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

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