

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092316\
 Data File : BG024102.D
 Acq On : 24 Sep 2016 00:31
 Operator : UM/SJ
 Sample : H4818-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-WATER2016-05-QT4

Manual Integrations
 APPROVED

sohil
 9/26/2016 6:41:10 PM

Quant Time: Sep 26 10:40:13 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG092316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 23 15:35:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	38080	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	169912	20.00	ng	0.00
38) Acenaphthene-d10	14.72	164	143549	20.00	ng	0.00
63) Phenanthrene-d10	17.48	188	372407	20.00	ng	0.00
75) Chrysene-d12	21.79	240	430633	20.00	ng	0.00
86) Perylene-d12	25.12	264	416706	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	314803	143.22	ng	0.00
7) Phenol-d6	7.21	99	479874	129.75	ng	0.00
23) Nitrobenzene-d5	9.22	82	348473	80.83	ng	0.00
41) 2,4,6-Tribromophenol	16.22	330	264693	128.28	ng	0.00
44) 2-Fluorobiphenyl	13.32	172	722394	84.59	ng	0.00
78) Terphenyl-d14	20.09	244	1516689	72.12	ng	0.00
Target Compounds						
32) Benzoic acid	10.34	122	2151m	3.25	ng	Qvalue # 83
40) Hexachlorocyclopentadiene	12.87	237	592m	3.98	ng	
52) 3-Nitroaniline	14.67	138	7279m	3.15	ng	
55) 4-Nitrophenol	15.08	139	1422	6.07	ng	
64) 4,6-Dinitro-2-methylphenol	15.90	198	5334	2.06	ng	
69) Pentachlorophenol	17.17	266	5108m	5.50	ng	

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

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