

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122118\
 Data File : BG038797.D
 Acq On : 21 Dec 2018 22:04
 Operator : JU/SJ
 Sample : J6503-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TP-GROUNDWATER-1-1

Quant Time: Dec 22 03:28:10 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	25583	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	109889	20.00	ng	0.00
39) Acenaphthene-d10	14.70	164	76144	20.00	ng	0.00
64) Phenanthrene-d10	17.45	188	179201	20.00	ng	0.00
76) Chrysene-d12	21.72	240	168097	20.00	ng	0.00
87) Perylene-d12	25.04	264	181813	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	103599	71.82	ng	0.01
7) Phenol-d6	7.22	99	89800	45.35	ng	0.01
23) Nitrobenzene-d5	9.24	82	199340	92.67	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	183035	174.42	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	552455	99.53	ng	0.00
79) Terphenyl-d14	20.05	244	871755	112.87	ng	0.00
Target Compounds						
10) Phenol	7.25	94	20016	9.515	ng	95
20) 3+4-Methylphenols	8.83	107	112211	57.648	ng	92
50) Dimethylphthalate	14.14	163	43301	6.964	ng	99

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

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