

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG122718\
 Data File : BG038849.D
 Acq On : 27 Dec 2018 18:40
 Operator : JU/SJ
 Sample : J6533-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 COMP-3

Quant Time: Dec 28 10:06:08 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	8.05	152	22757	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	95632	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	68396	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	167637	20.00	ng	0.00
76) Chrysene-d12	21.72	240	174308	20.00	ng	0.00
87) Perylene-d12	25.02	264	188470	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	109245	85.14	ng	0.00
7) Phenol-d6	7.24	99	101998	57.91	ng	0.03
23) Nitrobenzene-d5	9.24	82	180693	96.53	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	167479	177.67	ng	0.00
45) 2-Fluorobiphenyl	13.31	172	497001	99.69	ng	0.00
79) Terphenyl-d14	20.04	244	812245	101.41	ng	0.00
Target Compounds						
10) Phenol	7.26	94	12292	6.569	ng	84
15) Benzyl Alcohol	8.31	79	85693	62.330	ng	96
20) 3+4-Methylphenols	8.85	107	18710	10.806	ng	86
32) Benzoic acid	10.27	122	74015	76.288	ng	94

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

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