

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020719\
 Data File : BG039410.D
 Acq On : 8 Feb 2019 10:04
 Operator : JU/SJ
 Sample : K1401-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 RBR-200028

Quant Time: Feb 09 01:11:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG012919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 29 15:41:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.18	152	49124	20.00	ng	0.00
21) Naphthalene-d8	11.00	136	213441	20.00	ng	-0.01
39) Acenaphthene-d10	14.80	164	138773	20.00	ng	-0.01
64) Phenanthrene-d10	17.54	188	331561	20.00	ng	-0.01
76) Chrysene-d12	21.83	240	367830	20.00	ng	-0.02
87) Perylene-d12	25.22	264	349509	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	431989	150.51	ng	0.00
7) Phenol-d6	7.31	99	567710	134.37	ng	0.00
23) Nitrobenzene-d5	9.35	82	396682	116.10	ng	-0.01
42) 2,4,6-Tribromophenol	16.28	330	221222	148.04	ng	0.00
45) 2-Fluorobiphenyl	13.42	172	916160	94.71	ng	-0.01
79) Terphenyl-d14	20.14	244	1547776	89.07	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	11.05	128	188805	17.831	ng	99
37) 2-Methylnaphthalene	12.64	142	23206	2.959	ng	89
38) 1-Methylnaphthalene	12.86	142	16970	2.245	ng	91
52) Acenaphthene	14.86	154	27636	3.248	ng	97
55) Dibenzofuran	15.20	168	30054	2.203	ng	96
58) Fluorene	15.84	166	26757	2.631	ng	94
71) Phenanthrene	17.59	178	56307	3.281	ng	99
73) Carbazole	17.94	167	67110	3.978	ng	97

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

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