

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101717\
 Data File : BG029333.D
 Acq On : 18 Oct 2017 14:29
 Operator : SJ/JU
 Sample : I5906-14DL 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WH-TWS-04DL

Quant Time: Oct 18 16:39:38 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	66926	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	327345	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	215021	20.00	ng	0.00
63) Phenanthrene-d10	17.52	188	468212	20.00	ng	-0.01
75) Chrysene-d12	21.79	240	506440	20.00	ng	-0.01
86) Perylene-d12	25.10	264	510168	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.72	112	113393	28.30	ng	0.00
7) Phenol-d6	7.32	99	147386	26.21	ng	0.00
23) Nitrobenzene-d5	9.33	82	103388	20.07	ng	-0.01
41) 2,4,6-Tribromophenol	16.26	330	74164	26.71	ng	-0.01
44) 2-Fluorobiphenyl	13.41	172	289340	19.74	ng	0.00
78) Terphenyl-d14	20.11	244	450741	20.25	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	7.35	94	120962	19.952	ng	93
17) 2-Methylphenol	8.61	107	61385	15.242	ng	93
20) 3+4-Methylphenols	8.94	107	231568	40.808	ng	87
27) 2,4-Dimethylphenol	10.15	122	139025	27.759	ng	89
31) Naphthalene	11.04	128	1448062	88.761	ng	98
32) Benzoic acid	10.22	122	12027	2.868	ng	# 77
37) 2-Methylnaphthalene	12.63	142	139872	11.764	ng	92
51) Acenaphthene	14.85	154	103195	7.518	ng	97
54) Dibenzofuran	15.18	168	162496	8.292	ng	99
57) Fluorene	15.83	166	89005	5.498	ng	98
70) Phenanthrene	17.57	178	288628	11.933	ng	100
72) Carbazole	17.92	167	292566	12.540	ng	97
74) Fluoranthene	19.56	202	107585	3.803	ng	99
77) Pyrene	19.92	202	76524	2.491	ng	99

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

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