

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG012918\
 Data File : BG032420.D
 Acq On : 29 Jan 2018 18:41
 Operator : SJ/JU
 Sample : J1196-22DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 1035DL

Quant Time: Jan 30 04:26:44 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 29 15:32:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.72	152	24822	20.00	ng	0.00
21) Naphthalene-d8	11.60	136	100894	20.00	ng	0.00
38) Acenaphthene-d10	15.36	164	76739	20.00	ng	0.00
63) Phenanthrene-d10	18.09	188	216358	20.00	ng	0.00
75) Chrysene-d12	22.42	240	266081	20.00	ng	0.00
86) Perylene-d12	26.10	264	279367	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	6.15	112	37496	30.34	ng	0.00
7) Phenol-d6	7.83	99	48392	29.34	ng	0.00
23) Nitrobenzene-d5	9.92	82	30226	18.71	ng	0.00
41) 2,4,6-Tribromophenol	16.83	330	47012	32.18	ng	0.00
44) 2-Fluorobiphenyl	13.98	172	128950	19.97	ng	0.00
78) Terphenyl-d14	20.62	244	314159	25.29	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) 2-Chlorophenol	8.27	128	18457	12.635	ng	# 73
17) 2-Methylphenol	9.15	107	18331	14.412	ng	95
20) 3+4-Methylphenols	9.48	107	49694	27.853	ng	96
26) 2-Nitrophenol	10.69	139	18893	20.009	ng	# 85
27) 2,4-Dimethylphenol	10.73	122	25958	19.154	ng	94
29) 2,4-Dichlorophenol	11.23	162	36872	20.538	ng	87
36) 4-Chloro-3-methylphenol	12.81	107	60401	35.468	ng	# 93
42) 2,4,6-Trichlorophenol	13.80	196	66614	34.112	ng	95
43) 2,4,5-Trichlorophenol	13.87	196	20085	8.823	ng	# 91
53) 2,4-Dinitrophenol	15.45	184	25180	29.873	ng	# 54
55) 4-Nitrophenol	15.52	139	34289	35.308	ng	# 78
64) 4,6-Dinitro-2-methylphenol	16.45	198	27547	17.043	ng	# 72

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

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