

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
 Data File : BF085803.D
 Acq On : 28 Mar 2016 16:00
 Operator : UM/SJ
 Sample : H1739-06DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WPDL:

Quant Time: Mar 28 17:14:39 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 28 15:50:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	112060	20.00	ng	-0.01
21) Naphthalene-d8	8.10	136	454975	20.00	ng	-0.01
38) Acenaphthene-d10	9.85	164	206401	20.00	ng	-0.01
63) Phenanthrene-d10	11.34	188	407472	20.00	ng	-0.01
75) Chrysene-d12	13.97	240	281405	20.00	ng	-0.01
86) Perylene-d12	15.39	264	195487	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	184757	27.46	ng	0.00
7) Phenol-d6	6.45	99	249685	29.19	ng	-0.01
23) Nitrobenzene-d5	7.37	82	140424	17.38	ng	-0.02
41) 2,4,6-Tribromophenol	10.64	330	56230	28.67	ng	-0.01
44) 2-Fluorobiphenyl	9.18	172	302897	24.52	ng	-0.01
78) Terphenyl-d14	12.92	244	256884	18.77	ng	-0.01
Target Compounds						Qvalue
8) 2-Chlorophenol	6.60	128	223813	28.63	ng	97
17) 2-Methylphenol	7.08	107	159946	25.56	ng	96
20) 3+4-Methylphenols	7.22	107	255334	33.93	ng	# 55
26) 2-Nitrophenol	7.72	139	77072	18.50	ng	# 81
27) 2,4-Dimethylphenol	7.76	122	141059	19.37	ng	94
29) 2,4-Dichlorophenol	7.97	162	129991	21.23	ng	92
36) 4-Chloro-3-methylphenol	8.65	107	108129	15.12	ng	94
42) 2,4,6-Trichlorophenol	9.10	196	119702	28.96	ng	97
43) 2,4,5-Trichlorophenol	9.13	196	77534	17.88	ng	# 87
53) 2,4-Dinitrophenol	9.92	184	24452	19.31	ng	94
55) 4-Nitrophenol	9.98	139	70942	28.22	ng	96
64) 4,6-Dinitro-2-methylphenol	10.46	198	18347	7.69	ng	# 65
69) Pentachlorophenol	11.14	266	54338	25.39	ng	98

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

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