

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031518\
 Data File : BF103678.D
 Acq On : 15 Mar 2018 19:48
 Operator : JU/SJ
 Sample : J1888-06DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 16 04:55:56 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	153852	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	647885	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	308728	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	564855	20.00	ng	0.00
75) Chrysene-d12	14.14	240	483669	20.00	ng	0.00
86) Perylene-d12	15.67	264	447640	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	288355	28.51	ng	0.00
7) Phenol-d6	6.57	99	354324	28.63	ng	-0.01
23) Nitrobenzene-d5	7.52	82	180692	19.45	ng	-0.01
41) 2,4,6-Tribromophenol	10.79	330	80715	30.41	ng	0.00
44) 2-Fluorobiphenyl	9.32	172	418726	21.63	ng	0.00
78) Terphenyl-d14	13.08	244	464163	22.28	ng	0.00
Target Compounds						
8) 2-Chlorophenol	6.75	128	306448	28.920	ng	98
10) Phenol	6.59	94	281826	21.432	ng	100
17) 2-Methylphenol	7.20	107	92034	9.753	ng	98
20) 3+4-Methylphenols	7.35	107	205710	17.178	ng	87
26) 2-Nitrophenol	7.86	139	78266	24.595	ng	97
27) 2,4-Dimethylphenol	7.89	122	235487	25.559	ng	97
29) 2,4-Dichlorophenol	8.09	162	266430	31.784	ng	99
36) 4-Chloro-3-methylphenol	8.77	107	278453	30.132	ng	99
42) 2,4,6-Trichlorophenol	9.23	196	189161	33.577	ng	99
43) 2,4,5-Trichlorophenol	9.26	196	58692	9.230	ng	98
55) 4-Nitrophenol	10.09	139	88584	24.235	ng	94
64) 4,6-Dinitro-2-methylphenol	10.59	198	22221	18.996	ng	90
69) Pentachlorophenol	11.29	266	54703	14.375	ng	96

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

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