

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031518\
 Data File : BF103677.D
 Acq On : 15 Mar 2018 19:21
 Operator : JU/SJ
 Sample : J1888-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Mar 16 04:51:16 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	6.97	152	163016	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	678001	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	319161	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	584413	20.00	ng	0.00
75) Chrysene-d12	14.14	240	506360	20.00	ng	0.00
86) Perylene-d12	15.67	264	464479	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.59	112	1208351	112.74	ng	0.01
7) Phenol-d6	6.59	99	1515910	115.61	ng	0.00
23) Nitrobenzene-d5	7.53	82	886376	91.17	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	365912	133.34	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1563490	78.14	ng	0.00
78) Terphenyl-d14	13.09	244	1672348	76.66	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
8) 2-Chlorophenol	6.76	128	1201229	106.990	ng	97
10) Phenol	6.61	94	1178640	84.591	ng	96
17) 2-Methylphenol	7.20	107	419504	41.958	ng	99
20) 3+4-Methylphenols	7.36	107	900768	70.990	ng	87
26) 2-Nitrophenol	7.87	139	440906	90.995	ng	95
27) 2,4-Dimethylphenol	7.89	122	1004179	104.148	ng	98
29) 2,4-Dichlorophenol	8.10	162	1048438	119.520	ng	98
36) 4-Chloro-3-methylphenol	8.79	107	1094474	113.176	ng	98
42) 2,4,6-Trichlorophenol	9.24	196	757207	130.015	ng	98
43) 2,4,5-Trichlorophenol	9.27	196	245116	37.289	ng	99
55) 4-Nitrophenol	10.11	139	501538	110.481	ng	95
64) 4,6-Dinitro-2-methylphenol	10.60	198	201786	74.858	ng	83
69) Pentachlorophenol	11.29	266	267685	67.987	ng	97

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

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