

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031815\
 Data File : BF077801.D
 Acq On : 18 Mar 2015 17:46
 Operator : TP/IZ
 Sample : G1457-06DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 19 01:49:27 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 13 14:04:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.16	152	40788	20.00	ng	-0.01
21) Naphthalene-d8	8.74	136	169864	20.00	ng	-0.01
38) Acenaphthene-d10	10.91	164	82784	20.00	ng	0.00
63) Phenanthrene-d10	12.74	188	170488	20.00	ng	-0.01
75) Chrysene-d12	16.03	240	175668	20.00	ng	0.00
86) Perylene-d12	17.76	264	170455	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	86706	37.11	ng	-0.02
7) Phenol-d6	6.73	99	114667	39.72	ng	-0.01
23) Nitrobenzene-d5	7.85	82	61659	23.79	ng	-0.01
41) 2,4,6-Tribromophenol	11.88	330	28588	36.09	ng	-0.01
44) 2-Fluorobiphenyl	10.08	172	153819	27.31	ng	-0.01
78) Terphenyl-d14	14.74	244	185828	25.84	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	6.89	128	31159	11.20	ng	96
10) Phenol	6.74	94	82183	24.08	ng	# 70
17) 2-Methylphenol	7.49	107	47147	20.82	ng	98
20) 3+4-Methylphenols	7.69	107	37273	12.54	ng	85
26) 2-Nitrophenol	8.27	139	41497	30.36	ng	92
27) 2,4-Dimethylphenol	8.34	122	25286	10.16	ng	98
29) 2,4-Dichlorophenol	8.57	162	31042	13.08	ng	94
36) 4-Chloro-3-methylphenol	9.45	107	56236	23.55	ng	96
42) 2,4,6-Trichlorophenol	9.97	196	34820	20.36	ng	98
43) 2,4,5-Trichlorophenol	10.02	196	63942	34.60	ng	97
53) 2,4-Dinitrophenol	11.00	184	5278	18.21	ng	# 77
55) 4-Nitrophenol	11.08	139	28668	26.18	ng	# 79
64) 4,6-Dinitro-2-methylphenol	11.65	198	13235	19.27	ng	86
69) Pentachlorophenol	12.51	266	32550	32.97	ng	97

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

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