

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032416\
 Data File : BF085752.D
 Acq On : 25 Mar 2016 9:54
 Operator : UM/SJ
 Sample : H1739-06DL 5X
 Misc :
 ALS Vial : 98 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 25 14:31:02 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 24 14:15:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	100867	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	380956	20.00	ng	-0.01
38) Acenaphthene-d10	9.86	164	166707	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	253430	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	195149	20.00	ng	0.00
86) Perylene-d12	15.40	264	191466	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	196877	32.51	ng	0.01
7) Phenol-d6	6.46	99	247278	32.11	ng	0.00
23) Nitrobenzene-d5	7.38	82	152347	22.52	ng	-0.01
41) 2,4,6-Tribromophenol	10.65	330	41271	26.06	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	273276	27.39	ng	0.00
78) Terphenyl-d14	12.93	244	177667	18.72	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	6.61	128	239464	34.03	ng	98
17) 2-Methylphenol	7.08	107	160854	28.55	ng	97
20) 3+4-Methylphenols	7.24	107	281322	41.53	ng	# 53
26) 2-Nitrophenol	7.73	139	79714	22.85	ng	95
27) 2,4-Dimethylphenol	7.76	122	137181	22.50	ng	97
29) 2,4-Dichlorophenol	7.97	162	141913	27.68	ng	97
36) 4-Chloro-3-methylphenol	8.66	107	99926	16.69	ng	93
42) 2,4,6-Trichlorophenol	9.10	196	101871	30.51	ng	98
43) 2,4,5-Trichlorophenol	9.14	196	66676	19.04	ng	# 94
53) 2,4-Dinitrophenol	9.94	184	1958	7.35	ng	83
55) 4-Nitrophenol	9.99	139	51406	25.32	ng	97
69) Pentachlorophenol	11.16	266	35444	26.63	ng	98

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

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