

Data Path : Z:\HPCHEM1\BNA M\DATA\BM072617\
 Data File : BM010972.D
 Acq On : 26 Jul 2017 16:00
 Operator : SJ/JU
 Sample : I4351-22
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 0535

Quant Time: Jul 26 17:02:17 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM072617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 26 17:01:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	253924	20.00	ng	0.00
21) Naphthalene-d8	10.15	136	1037235	20.00	ng	0.00
38) Acenaphthene-d10	14.05	164	705368	20.00	ng	0.00
63) Phenanthrene-d10	16.81	188	1716839	20.00	ng	0.00
75) Chrysene-d12	21.03	240	1820543	20.00	ng	0.00
86) Perylene-d12	23.14	264	1318822	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.03	112	2151615	143.24	ng	0.00
7) Phenol-d6	6.60	99	3041434	142.51	ng	0.00
23) Nitrobenzene-d5	8.55	82	2340899	84.70	ng	0.00
41) 2,4,6-Tribromophenol	15.56	330	1618797	143.19	ng	0.00
44) 2-Fluorobiphenyl	12.66	172	4967722	88.33	ng	0.00
78) Terphenyl-d14	19.47	244	8088607	88.01	ng	0.00

Target Compounds				Qvalue		
8) 2-Chlorophenol	6.97	128	2733992	179.273	ng	85
10) Phenol	6.63	94	2507825	108.184	ng	97
17) 2-Methylphenol	7.85	107	2040466	137.726	ng	# 81
20) 3+4-Methylphenols	8.17	107	3051683	148.179	ng	85
27) 2,4-Dimethylphenol	9.36	122	831354	51.827	ng	# 73
29) 2,4-Dichlorophenol	9.81	162	822434	45.627	ng	91
36) 4-Chloro-3-methylphenol	11.46	107	1466382	63.003	ng	# 75
42) 2,4,6-Trichlorophenol	12.47	196	1988193	108.392	ng	99
43) 2,4,5-Trichlorophenol	12.53	196	3341592	176.880	ng	# 87
55) 4-Nitrophenol	14.30	139	862452	92.707	ng	98
64) 4,6-Dinitro-2-methylphenol	15.20	198	666480	57.343	ng	87
69) Pentachlorophenol	16.46	266	1374669	86.820	ng	98

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

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