

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN072618\
 Data File : BN002168.D
 Acq On : 26 Jul 2018 10:43
 Operator : JU/SJ
 Sample : J4086-22
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 1535

Quant Time: Jul 27 01:33:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 26 18:27:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	209762	20.00	ng	0.00
21) Naphthalene-d8	10.46	136	800842	20.00	ng	0.00
38) Acenaphthene-d10	14.32	164	409733	20.00	ng	0.00
63) Phenanthrene-d10	17.07	188	824108	20.00	ng	0.00
75) Chrysene-d12	21.26	240	736041	20.00	ng	-0.01
86) Perylene-d12	23.49	264	759154	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	1855077	142.38	ng	0.00
7) Phenol-d6	6.88	99	2466173	134.88	ng	0.00
23) Nitrobenzene-d5	8.84	82	1531601	98.08	ng	0.00
41) 2,4,6-Tribromophenol	15.82	330	758662	142.59	ng	0.00
44) 2-Fluorobiphenyl	12.94	172	3161724	100.51	ng	0.00
78) Terphenyl-d14	19.71	244	3749517	106.25	ng	0.00
Target Compounds						
10) Phenol	6.91	94	2914856	155.524	ng	Qvalue 97
17) 2-Methylphenol	8.13	107	1405076	109.812	ng	95
20) 3+4-Methylphenols	8.46	107	1866754	104.567	ng	95
26) 2-Nitrophenol	9.59	139	1027858	131.475	ng	95
29) 2,4-Dichlorophenol	10.12	162	1634024	128.880	ng	95
36) 4-Chloro-3-methylphenol	11.76	107	1616785	112.146	ng	95
42) 2,4,6-Trichlorophenol	12.75	196	385504	41.704	ng	97
43) 2,4,5-Trichlorophenol	12.82	196	1188897	121.419	ng	96
53) 2,4-Dinitrophenol	14.45	184	656683	164.016	ng	98
55) 4-Nitrophenol	14.56	139	960820	152.757	ng	87
64) 4,6-Dinitro-2-methylphenol	15.46	198	220080	39.555	ng	98
69) Pentachlorophenol	16.72	266	790939	115.148	ng	99

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

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