

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061915\
 Data File : BF080062.D
 Acq On : 19 Jun 2015 17:29
 Operator : TP/IZ
 Sample : G2653-05
 Misc : LOD-SVOC-WATER-8PPM
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-WATER2015-01

Quant Time: Jun 20 00:49:05 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 17 16:41:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	39102	20.00	ng	-0.01
21) Naphthalene-d8	8.62	136	161276	20.00	ng	-0.01
38) Acenaphthene-d10	10.39	164	86137	20.00	ng	-0.01
63) Phenanthrene-d10	11.90	188	179502	20.00	ng	-0.01
75) Chrysene-d12	14.93	240	212437	20.00	ng	0.07
86) Perylene-d12	16.84	264	197819	20.00	ng	0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.94	112	238803	108.11	ng	0.00
7) Phenol-d6	6.93	99	307238	104.52	ng	-0.01
23) Nitrobenzene-d5	7.89	82	185499	66.96	ng	-0.01
41) 2,4,6-Tribromophenol	11.19	330	90937	107.96	ng	-0.01
44) 2-Fluorobiphenyl	9.69	172	375826	62.22	ng	-0.01
78) Terphenyl-d14	13.63	244	532667	58.56	ng	0.04
Target Compounds						Qvalue
49) Dimethylphthalate	10.08	163	31895	4.93	ng	# 97
59) Diethylphthalate	10.79	149	34841	5.34	ng	99
64) 4,6-Dinitro-2-methylphenol	10.97	198	1985	6.23	ng	87
66) 4-Bromophenyl-phenylether	11.43	248	9577	5.02	ng	# 92
68) Atrazine	11.56	200	8087	4.38	ng	84
70) Phenanthrene	11.92	178	52810	4.86	ng	97
72) Carbazole	12.13	167	43833	4.39	ng	95
74) Fluoranthene	13.20	202	52122	4.50	ng	94
77) Pyrene	13.47	202	56235	4.30	ng	97
80) Benzo(a)anthracene	14.92	228	56907	4.40	ng	97
82) Chrysene	14.96	228	54044	4.53	ng	95
84) Di-n-octyl phthalate	15.73	149	40567	2.85	ng	# 98
87) Benzo(b)fluoranthene	16.29	252	53525	4.47	ng	# 90
89) Benzo(a)pyrene	16.76	252	50164	4.41	ng	# 85
91) Benzo(g,h,i)perylene	19.25	276	49940	4.25	ng	# 76

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

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