

Data Path : Z:\HPCHEM1\BNA E\DATA\BE121417\
 Data File : BE094991.D
 Acq On : 14 Dec 2017 23:44
 Operator : SJ/JU
 Sample : I6797-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 YT-MW-01

Quant Time: Dec 15 08:31:46 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 14 18:42:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.47	152	7486	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	16916	0.40	ng	0.00
13) Acenaphthene-d10	15.06	164	9216	0.40	ng	0.00
19) Phenanthrene-d10	17.77	188	20611	0.40	ng	0.00
27) Chrysene-d12	21.87	240	25246	0.40	ng	0.00
34) Perylene-d12	24.54	264	19227	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.95	112	1783	0.15	ng	0.00
5) Phenol-d6	7.59	99	1678	0.10	ng	0.00
8) Nitrobenzene-d5	9.65	82	5355	0.39	ng	0.00
11) 2-Methylnaphthalene-d10	12.86	152	8943	0.32	ng	0.00
14) 2,4,6-Tribromophenol	16.53	330	1056	0.49	ng	0.00
15) 2-Fluorobiphenyl	13.70	172	16487	0.40	ng	0.00
25) Fluoranthene-d10	19.75	212	95243	0.36	ng	0.00
29) Terphenyl-d14	20.31	244	21794	0.46	ng	0.00
Target Compounds						Qvalue
22) Pentachlorophenol	17.41	266	149	0.128	ng	93
28) Pyrene	20.13	202	3799	0.044	ng	97
33) Indeno(1,2,3-cd)pyrene	27.40	276	2142	0.032	ng	95
35) Benzo(b)fluoranthene	23.71	252	2507	0.036	ng	# 58
38) Dibenzo(a,h)anthracene	27.43	278	1652	0.028	ng	# 50
39) Benzo(g,h,i)perylene	28.29	276	1878	0.032	ng	# 61

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

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