

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091117\
 Data File : BF098441.D
 Acq On : 12 Sep 2017 9:46
 Operator : SJ/JU
 Sample : I5151-03DL 250X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-PH4-ESMI2-3DL

Manual Integrations
 APPROVED

Sohil
 9/12/2017 6:02:29 PM

Quant Time: Sep 12 12:02:59 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	230587	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	933019	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	378735	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	548716	20.00	ng	0.00
75) Chrysene-d12	13.82	240	375183	20.00	ng	0.00
86) Perylene-d12	15.22	264	380917	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	4882	0.31	ng	0.02
7) Phenol-d6	6.34	99	8062	0.41	ng	0.02
23) Nitrobenzene-d5	7.24	82	5099	0.30	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	418	0.17	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	9371	0.42	ng	-0.01
78) Terphenyl-d14	12.76	244	5088	0.29	ng	0.00

Target Compounds

						Qvalue
31) Naphthalene	7.98	128	2890824	62.675	ng	99
37) 2-Methylnaphthalene	8.66	142	606207	21.697	ng	98
45) 1,1'-Biphenyl	9.13	154	186810	5.511	ng	98
48) Acenaphthylene	9.56	152	115212	3.201	ng	# 95
51) Acenaphthene	9.73	154	644222	28.155	ng	98
54) Dibenzofuran	9.90	168	542833	18.008	ng	98
57) Fluorene	10.25	166	463156	18.563	ng	99
70) Phenanthrene	11.21	178	1856753	63.830	ng	99
71) Anthracene	11.26	178	567296	18.995	ng	98
72) Carbazole	11.42	167	162771	5.848	ng	97
74) Fluoranthene	12.39	202	1119864	38.456	ng	98
77) Pylene	12.62	202	951401	32.145	ng	99
80) Benzo(a)anthracene	13.80	228	356470m	16.206	ng	
82) Chrysene	13.84	228	288343m	12.968	ng	
85) Indeno(1,2,3-cd)pyrene	16.55	276	132141	8.471	ng	96
87) Benzo(b)fluoranthene	14.82	252	387558m	16.181	ng	
88) Benzo(k)fluoranthene	14.84	252	106825m	4.778	ng	
89) Benzo(a)pyrene	15.16	252	304893	14.289	ng	96
91) Benzo(a,h,i)perylene	16.96	276	121745	7.798	ng	99

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

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