

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
 Data File : BF098397.D
 Acq On : 9 Sep 2017 20:30
 Operator : SJ/JU
 Sample : I5151-03 50X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 9/11/2017 2:42:52 PM

Quant Time: Sep 11 12:12:40 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	99381	20.00	ng	-0.03
21) Naphthalene-d8	7.96	136	368572	20.00	ng	-0.02
38) Acenaphthene-d10	9.71	164	169402	20.00	ng	-0.03
63) Phenanthrene-d10	11.19	188	268934	20.00	ng	-0.03
75) Chrysene-d12	13.83	240	192303	20.00	ng	-0.04
86) Perylene-d12	15.23	264	196243	20.00	ng	-0.03

System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	14611	2.24	ng	-0.02
7) Phenol-d6	6.33	99	22082	2.49	ng	-0.02
23) Nitrobenzene-d5	7.24	82	12426	1.83	ng	-0.04
41) 2,4,6-Tribromophenol	10.50	330	2657	1.81	ng	-0.04
44) 2-Fluorobiphenyl	9.03	172	22205	1.93	ng	-0.04
78) Terphenyl-d14	12.77	244	15113	1.64	ng	-0.04

Target Compounds				Qvalue		
31) Naphthalene	8.00	128	3986350	208.334	ng	99
37) 2-Methylnaphthalene	8.67	142	1048042	89.338	ng	99
45) 1,1'-Biphenyl	9.13	154	375109	24.282	ng	97
48) Acenaphthylene	9.57	152	278687	15.548	ng	# 96
51) Acenaphthene	9.75	154	1056832	93.488	ng	99
54) Dibenzofuran	9.92	168	1056058	69.704	ng	99
57) Fluorene	10.26	166	892621	76.204	ng	98
70) Phenanthrene	11.23	178	2955016	206.201	ng	99
71) Anthracene	11.27	178	1118329	76.939	ng	100
72) Carbazole	11.43	167	390701	28.318	ng	97
74) Fluoranthene	12.42	202	2155012	144.072	ng	100
77) Pvrene	12.65	202	1894811	120.472	ng	98
80) Benzo(a)anthracene	13.82	228	836050m	72.539	ng	
82) Chrysene	13.86	228	670086m	58.445	ng	
85) Indeno(1,2,3-cd)pyrene	16.58	276	337788m	40.050	ng	
87) Benzo(b)fluoranthene	14.84	252	892097m	72.119	ng	
88) Benzo(k)fluoranthene	14.86	252	272641m	23.460	ng	
89) Benzo(a)pyrene	15.18	252	702903	64.188	ng	97
90) Dibenzo(a,h)anthracene	16.59	278	69360m	7.780	ng	
91) Benzo(a,h,i)perylene	16.99	276	310423	33.371	ng	94

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

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