

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\
 Data File : BF098395.D
 Acq On : 9 Sep 2017 19:35
 Operator : SJ/JU
 Sample : I5151-01 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-PH4-ESMI2-1

Manual Integrations
 APPROVED

Sohil
 9/11/2017 2:42:48 PM

Quant Time: Sep 11 06:02:49 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	98908	20.00	ng	-0.03
21) Naphthalene-d8	7.96	136	377052	20.00	ng	-0.03
38) Acenaphthene-d10	9.70	164	160730	20.00	ng	-0.04
63) Phenanthrene-d10	11.19	188	267201	20.00	ng	-0.04
75) Chrysene-d12	13.82	240	223688	20.00	ng	-0.04
86) Perylene-d12	15.23	264	208767	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	159730	24.56	ng	-0.02
7) Phenol-d6	6.32	99	217708	24.64	ng	-0.03
23) Nitrobenzene-d5	7.24	82	123318	17.77	ng	-0.04
41) 2,4,6-Tribromophenol	10.50	330	28950	20.76	ng	-0.04
44) 2-Fluorobiphenyl	9.03	172	215179	19.67	ng	-0.04
78) Terphenyl-d14	12.77	244	141186	13.16	ng	-0.04
Target Compounds						Qvalue
31) Naphthalene	7.98	128	322122	16.456	ng	99
37) 2-Methylnaphthalene	8.67	142	209805	17.482	ng	99
45) 1,1'-Biphenyl	9.13	154	56221	3.836	ng	96
51) Acenaphthene	9.74	154	206859	19.286	ng	99
54) Dibenzofuran	9.91	168	206827	14.388	ng	99
57) Fluorene	10.25	166	217777	19.595	ng	96
70) Phenanthrene	11.22	178	736107	51.699	ng	100
71) Anthracene	11.26	178	268037	18.560	ng	99
72) Carbazole	11.42	167	39044	2.848	ng	97
74) Fluoranthene	12.40	202	493524	33.208	ng	97
77) Pyrene	12.63	202	395486	21.617	ng	97
80) Benzo(a)anthracene	13.81	228	207529m	15.480	ng	
82) Chrysene	13.84	228	147143	11.033	ng	96
85) Indeno(1,2,3-cd)pyrene	16.56	276	50821m	5.180	ng	
87) Benzo(b)fluoranthene	14.83	252	176239m	13.393	ng	
88) Benzo(k)fluoranthene	14.85	252	50699m	4.101	ng	
89) Benzo(a)pyrene	15.16	252	131778	11.312	ng	98
91) Benzo(g,h,i)perylene	16.97	276	41615	4.205	ng	# 92

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

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF090917\
Data File : BF098395.D
Acq On : 9 Sep 2017 19:35
Operator : SJ/JU
Sample : I5151-01 5X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
NYSEG-PH4-ESM12-1

Manual Integrations
APPROVED

Sohil
9/11/2017 2:42:48 PM

Quant Time: Sep 11 06:02:49 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

