

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
 Data File : BF098978.D
 Acq On : 30 Sep 2017 22:26
 Operator : SJ/JU
 Sample : I5563-04 10X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-PH4-SB-ESMI-4

Manual Integrations
 APPROVED

Sohil
 10/2/2017 5:44:43 PM

Quant Time: Oct 01 01:27:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	130446	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	479135	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	177437	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	279186	20.00	ng	0.00
75) Chrysene-d12	14.09	240	229344	20.00	ng	0.00
86) Perylene-d12	15.59	264	166832	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	93190	11.37	ng	-0.01
7) Phenol-d6	6.53	99	116528	11.53	ng	-0.01
23) Nitrobenzene-d5	7.47	82	66492	8.90	ng	-0.01
41) 2,4,6-Tribromophenol	10.74	330	14118	9.72	ng	0.00
44) 2-Fluorobiphenyl	9.27	172	127406	11.99	ng	-0.01
78) Terphenyl-d14	13.02	244	95656	9.49	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.23	128	1810871	80.472	ng	99
37) 2-Methylnaphthalene	8.91	142	365205	24.965	ng	99
45) 1,1'-Biphenyl	9.37	154	145513	9.542	ng	96
48) Acenaphthylene	9.82	152	126333	7.208	ng	98
51) Acenaphthene	9.99	154	547195	49.284	ng	99
54) Dibenzofuran	10.16	168	473860	32.054	ng	99
57) Fluorene	10.50	166	459115	38.639	ng	99
70) Phenanthrene	11.48	178	2049667	137.156	ng	99
71) Anthracene	11.53	178	984937	64.415	ng	100
72) Carbazole	11.67	167	207492	13.857	ng	100
74) Fluoranthene	12.67	202	1915190	126.561	ng	100
77) Pylene	12.90	202	1675621	95.814	ng	99
80) Benzo(a)anthracene	14.08	228	778921m	56.089	ng	
82) Chrysene	14.12	228	585859	44.311	ng	98
85) Indeno(1,2,3-cd)pyrene	17.10	276	203375	19.229	ng	98
87) Benzo(b)fluoranthene	15.14	252	573565m	55.917	ng	
88) Benzo(k)fluoranthene	15.17	252	184046m	18.166	ng	
89) Benzo(a)pyrene	15.52	252	418164	44.412	ng	# 95
90) Dibenzo(a,h)anthracene	17.10	278	45434	6.029	ng	# 78
91) Benzo(a,h,i)perylene	17.56	276	191473	25.706	ng	98

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

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