

Data Path : Z:\HPCHEM1\BNA F\DATA\BF093017\
 Data File : BF098980.D
 Acq On : 30 Sep 2017 23:21
 Operator : SJ/JU
 Sample : I5563-07 10X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NYSEG-PH4-SB-ESMI-7

Manual Integrations
 APPROVED

Sohil
 10/2/2017 5:44:47 PM

Quant Time: Oct 01 01:37:29 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	120690	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	441754	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	160025	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	257961	20.00	ng	0.00
75) Chrysene-d12	14.09	240	215497	20.00	ng	0.00
86) Perylene-d12	15.59	264	157164	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	99579	13.13	ng	-0.01
7) Phenol-d6	6.53	99	118652	12.69	ng	-0.01
23) Nitrobenzene-d5	7.47	82	65929	9.57	ng	-0.02
41) 2,4,6-Tribromophenol	10.74	330	14359	10.97	ng	0.00
44) 2-Fluorobiphenyl	9.27	172	100773	10.51	ng	-0.01
78) Terphenyl-d14	13.02	244	75223	7.94	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.22	128	333915	16.094	ng	99
37) 2-Methylnaphthalene	8.91	142	42851	3.177	ng	97
45) 1,1'-Biphenyl	9.37	154	66963	4.869	ng	97
48) Acenaphthylene	9.82	152	117266	7.418	ng	98
51) Acenaphthene	9.99	154	308169	30.776	ng	99
54) Dibenzofuran	10.16	168	251862	18.891	ng	96
57) Fluorene	10.50	166	350535	32.711	ng	97
70) Phenanthrene	11.47	178	1768630	128.088	ng	99
71) Anthracene	11.53	178	910527	64.448	ng	99
72) Carbazole	11.67	167	126703	9.158	ng	100
74) Fluoranthene	12.67	202	2094538	149.802	ng	98
77) Pylene	12.90	202	1906730	116.034	ng	98
80) Benzo(a)anthracene	14.08	228	871379m	66.778	ng	
82) Chrysene	14.12	228	679428	54.691	ng	98
85) Indeno(1,2,3-cd)pyrene	17.10	276	270014	27.171	ng	96
87) Benzo(b)fluoranthene	15.14	252	657125m	68.004	ng	
88) Benzo(k)fluoranthene	15.17	252	254284m	26.643	ng	
89) Benzo(a)pyrene	15.53	252	523875	59.061	ng	96
90) Dibenzo(a,h)anthracene	17.11	278	55892	7.874	ng	# 81
91) Benzo(a,h,i)perylene	17.57	276	258555	36.848	ng	96

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

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