

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081117\
 Data File : BF097609.D
 Acq On : 11 Aug 2017 22:57
 Operator : SJ/JU
 Sample : I4697-08 50X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

sohil
 8/14/2017 6:54:41 PM

Quant Time: Aug 14 10:51:58 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 15:55:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	84270	20.00	ng	0.00
21) Naphthalene-d8	9.54	136	351282	20.00	ng	0.00
38) Acenaphthene-d10	12.37	164	141090	20.00	ng	0.00
63) Phenanthrene-d10	14.77	188	202381	20.00	ng	0.00
75) Chrysene-d12	18.47	240	154182	20.00	ng	0.00
86) Perylene-d12	20.13	264	175692	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.57	112	7083m	1.34	ng	0.17
7) Phenol-d6	7.19	99	12603m	1.92	ng	0.13
23) Nitrobenzene-d5	8.47	82	7105m	1.27	ng	0.05
41) 2,4,6-Tribromophenol	13.67	330	1213	1.13	ng	0.01
44) 2-Fluorobiphenyl	11.32	172	13166	1.41	ng	0.00
78) Terphenyl-d14	17.12	244	7670	1.03	ng	0.00

Target Compounds

						Qvalue
31) Naphthalene	9.59	128	2954377	167.909	ng	99
37) 2-Methylnaphthalene	10.70	142	611101	55.199	ng	98
45) 1,1'-Biphenyl	11.46	154	245589	20.289	ng	95
48) Acenaphthylene	12.13	152	176004	11.912	ng	98
51) Acenaphthene	12.43	154	853429	96.381	ng	98
54) Dibenzofuran	12.71	168	739543	59.659	ng	99
57) Fluorene	13.26	166	705895	69.243	ng	99
70) Phenanthrene	14.83	178	2591113	222.292	ng	100
71) Anthracene	14.90	178	810018	67.919	ng	99
72) Carbazole	15.18	167	250494	22.063	ng	98
74) Fluoranthene	16.58	202	1723329	163.615	ng	98
77) Pyrene	16.89	202	1434317	112.494	ng	100
80) Benzo(a)anthracene	18.46	228	666449	74.009	ng	94
82) Chrysene	18.50	228	466114	52.041	ng	96
85) Indeno(1,2,3-cd)pyrene	21.36	276	338598	48.606	ng	95
87) Benzo(b)fluoranthene	19.73	252	607391m	57.573	ng	
88) Benzo(k)fluoranthene	19.76	252	249888m	23.518	ng	
89) Benzo(a)pyrene	20.08	252	580551	59.953	ng	96
90) Dibenzo(a,h)anthracene	21.36	278	63386	9.142	ng	# 76
91) Benzo(a,h,i)perylene	21.71	276	326693	42.953	ng	97

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

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