

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081717\
 Data File : BF097820.D
 Acq On : 18 Aug 2017 7:53
 Operator : SJ/JU
 Sample : I4773-01DL 50X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-J10-ESWDL

Manual Integrations
 APPROVED

Sohil
 8/18/2017 7:02:39 PM

Quant Time: Aug 18 08:35:29 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 15 13:25:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	127168	20.00	ng	-0.01
21) Naphthalene-d8	8.06	136	459599	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	159072	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	271078	20.00	ng	0.00
75) Chrysene-d12	13.93	240	281861	20.00	ng	0.00
86) Perylene-d12	15.36	264	199396	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.37	112	17668	2.11	ng	-0.02
7) Phenol-d6	6.39	99	24092	2.12	ng	-0.02
23) Nitrobenzene-d5	7.33	82	13923	1.65	ng	-0.02
41) 2,4,6-Tribromophenol	10.60	330	2013	1.46	ng	-0.01
44) 2-Fluorobiphenyl	9.13	172	16385	1.51	ng	-0.01
78) Terphenyl-d14	12.88	244	12726	0.94	ng	-0.01

Target Compounds

						Qvalue
12) 1,3-Dichlorobenzene	6.72	146	35467	3.740	ng	# 93
13) 1,4-Dichlorobenzene	6.79	146	126840	13.150	ng	# 92
14) 1,2-Dichlorobenzene	6.95	146	90566	10.183	ng	97
30) 1,2,4-Trichlorobenzene	8.00	180	23461	3.475	ng	97
31) Naphthalene	8.09	128	1525821	63.949	ng	99
37) 2-Methylnaphthalene	8.77	142	907298	62.023	ng	98
45) 1,1'-Biphenyl	9.23	154	174358	12.020	ng	95
51) Acenaphthene	9.85	154	262836	24.760	ng	99
54) Dibenzofuran	10.02	168	216509	15.218	ng	# 85
57) Fluorene	10.36	166	225649	20.515	ng	95
70) Phenanthrene	11.32	178	732837	50.733	ng	100
71) Anthracene	11.37	178	139295	9.508	ng	99
72) Carbazole	11.52	167	28208	2.028	ng	# 90
74) Fluoranthene	12.51	202	367907	24.402	ng	98
77) Pyrene	12.73	202	319595	13.863	ng	98
80) Benzo(a)anthracene	13.92	228	83391	4.936	ng	94
82) Chrysene	13.96	228	72063	4.288	ng	97
87) Benzo(b)fluoranthene	14.94	252	54554m	4.341	ng	
89) Benzo(a)pyrene	15.30	252	29688	2.668	ng	98
91) Benzo(a,h,i)perylene	17.17	276	21073	2.230	ng	94

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

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