

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090817\
 Data File : BF098351.D
 Acq On : 8 Sep 2017 15:34
 Operator : SJ/JU
 Sample : I5095-01DL 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STORM-SEWER-SEDDL

Manual Integrations
 APPROVED

Sohil
 9/11/2017 4:47:34 PM

Quant Time: Sep 08 16:09:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 08 15:02:28 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	101480	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	398972	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	166973	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	275820	20.00	ng	0.00
75) Chrysene-d12	13.83	240	200991	20.00	ng	0.00
86) Perylene-d12	15.24	264	218669	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	298014	44.67	ng	0.00
7) Phenol-d6	6.32	99	398453	43.96	ng	0.00
23) Nitrobenzene-d5	7.25	82	236180	32.16	ng	0.00
41) 2,4,6-Tribromophenol	10.50	330	54704	37.76	ng	0.00
44) 2-Fluorobiphenyl	9.04	172	288765	25.41	ng	0.00
78) Terphenyl-d14	12.78	244	164262	17.04	ng	0.00

Target Compounds

						Qvalue
48) Acenaphthylene	9.57	152	48506	2.746	ng	99
49) Dimethylphthalate	9.44	163	58013	4.753	ng	97
57) Fluorene	10.26	166	39132	3.389	ng	92
70) Phenanthrene	11.22	178	410128	27.904	ng	99
71) Anthracene	11.27	178	115809	7.769	ng	98
72) Carbazole	11.43	167	39815	2.814	ng	97
74) Fluoranthene	12.41	202	748617	48.799	ng	97
77) Pyrene	12.64	202	633561	38.541	ng	99
79) Butylbenzylphthalate	13.26	149	13834	2.195	ng	# 65
80) Benzo(a)anthracene	13.83	228	368016	30.550	ng	96
82) Chrysene	13.86	228	352149	29.387	ng	98
83) Bis(2-ethylhexyl)phthalate	13.82	149	47989	6.871	ng	# 99
87) Benzo(b)fluoranthene	14.85	252	457415m	33.186	ng	
88) Benzo(k)fluoranthene	14.87	252	166317m	12.843	ng	
89) Benzo(a)pyrene	15.19	252	343169	28.124	ng	97
90) Dibenzo(a,h)anthracene	16.60	278	56863m	5.724	ng	
91) Benzo(a,h,i)perylene	17.01	276	190705	18.398	ng	94

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

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