

Data Path : Z:\HPCHEM1\BNA F\DATA\BF032917\
 Data File : BF094058.D
 Acq On : 30 Mar 2017 12:51
 Operator : SJ/MA
 Sample : I2443-11DL 5X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-AA9-SSWDL

Quant Time: Mar 30 13:20:29 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 27 16:10:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.92	152	154189	20.00	ng	-0.03
21) Naphthalene-d8	7.43	136	587563	20.00	ng	-0.03
38) Acenaphthene-d10	9.57	164	245427	20.00	ng	-0.03
63) Phenanthrene-d10	11.40	188	377729	20.00	ng	-0.02
75) Chrysene-d12	14.65	240	297518	20.00	ng	-0.03
86) Perylene-d12	16.29	264	220961	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	4.46	112	232922	25.51	ng	-0.01
7) Phenol-d6	5.53	99	278529	24.66	ng	-0.02
23) Nitrobenzene-d5	6.56	82	102457	9.95	ng	-0.04
41) 2,4,6-Tribromophenol	10.55	330	34499	17.79	ng	-0.03
44) 2-Fluorobiphenyl	8.75	172	282451	17.55	ng	-0.03
78) Terphenyl-d14	13.37	244	180687	13.61	ng	-0.03
Target Compounds						Qvalue
31) Naphthalene	7.45	128	96109	3.40	ng	# 95
51) Acenaphthene	9.61	154	65162	4.34	ng	95
54) Dibenzofuran	9.82	168	53611	2.62	ng	95
57) Fluorene	10.25	166	92806	5.47	ng	99
70) Phenanthrene	11.43	178	768993	36.60	ng	98
71) Anthracene	11.49	178	220412	10.19	ng	98
72) Carbazole	11.69	167	69207	3.12	ng	97
74) Fluoranthene	12.89	202	713445	33.16	ng	99
77) Pyrene	13.17	202	626361	29.01	ng	98
80) Benzo(a)anthracene	14.64	228	269975m	15.56	ng	
82) Chrysene	14.69	228	254486m	15.81	ng	
85) Indeno(1,2,3-cd)pyrene	17.65	276	72398	5.19	ng	# 91
87) Benzo(b)fluoranthene	15.88	252	216224m	15.96	ng	
88) Benzo(k)fluoranthene	15.90	252	79901m	6.03	ng	
89) Benzo(a)pyrene	16.23	252	162764	13.05	ng	98
91) Benzo(g,h,i)perylene	18.05	276	71629	6.73	ng	96

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

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