

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011017\
 Data File : BF092196.D
 Acq On : 10 Jan 2017 21:44
 Operator : UM/SJ
 Sample : I1098-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP3-COMP13

Quant Time: Jan 11 00:58:00 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 06 15:43:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.62	152	145184	20.00	ng	-0.02
21) Naphthalene-d8	7.91	136	630791	20.00	ng	-0.02
38) Acenaphthene-d10	9.67	164	343425	20.00	ng	-0.01
63) Phenanthrene-d10	11.14	188	583909	20.00	ng	-0.01
75) Chrysene-d12	13.79	240	390291	20.00	ng	0.00
86) Perylene-d12	15.16	264	313995	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.24	112	1480434	170.26	ng	0.00
7) Phenol-d6	6.30	99	1617841	152.40	ng	-0.01
23) Nitrobenzene-d5	7.20	82	1280442	112.87	ng	-0.01
41) 2,4,6-Tribromophenol	10.46	330	444821	156.51	ng	-0.01
44) 2-Fluorobiphenyl	9.00	172	1778525	109.89	ng	-0.01
78) Terphenyl-d14	12.73	244	1778443	110.51	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.31	94	119182	9.85	ng	82
49) Dimethylphthalate	9.40	163	374561	17.05	ng	100
70) Phenanthrene	11.17	178	314275	9.93	ng	97
74) Fluoranthene	12.36	202	525982	16.51	ng	93
77) Pyrene	12.58	202	501219	17.50	ng	97
80) Benzo(a)anthracene	13.77	228	281569	12.78	ng	97
82) Chrysene	13.81	228	230821m	10.44	ng	
85) Indeno(1,2,3-cd)pyrene	16.43	276	179806	9.39	ng	94
87) Benzo(b)fluoranthene	14.78	252	326296m	16.69	ng	
88) Benzo(k)fluoranthene	14.79	252	99668m	5.98	ng	
89) Benzo(a)pyrene	15.10	252	277437	15.99	ng	97
90) Dibenzo(a,h)anthracene	16.43	278	65731	4.61	ng	# 83
91) Benzo(a,h,i)perylene	16.80	276	181765	12.26	ng	# 95

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

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