

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121816\
 Data File : BF091767.D
 Acq On : 19 Dec 2016 6:14
 Operator : UM/SJ
 Sample : H6006-22DL 5X
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-2-0-5DL

Quant Time: Dec 19 17:07:32 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 17 22:53:54 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	204280m	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	856407	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	395633	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	609630	20.00	ng	0.00
75) Chrysene-d12	13.92	240	449483	20.00	ng	-0.01
86) Perylene-d12	15.32	264	463488	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	111132	9.11	ng	-0.01
7) Phenol-d6	6.41	99	226983	15.26	ng	-0.01
23) Nitrobenzene-d5	7.32	82	123969	8.35	ng	-0.02
41) 2,4,6-Tribromophenol	10.59	330	56122	16.73	ng	-0.01
44) 2-Fluorobiphenyl	9.13	172	330620	16.79	ng	-0.01
78) Terphenyl-d14	12.87	244	278576	15.65	ng	-0.01
Target Compounds						Qvalue
70) Phenanthrene	11.31	178	620554	20.37	ng	97
71) Anthracene	11.36	178	136129	4.22	ng	98
72) Carbazole	11.52	167	72780	2.58	ng	98
74) Fluoranthene	12.50	202	875051	27.50	ng	91
77) Pyrene	12.72	202	679520	20.93	ng	98
80) Benzo(a)anthracene	13.91	228	398288	15.49	ng	98
82) Chrysene	13.94	228	290335	10.89	ng	97
85) Indeno(1,2,3-cd)pyrene	16.66	276	205608	7.85	ng	96
87) Benzo(b)fluoranthene	14.92	252	502772m	17.11	ng	
88) Benzo(k)fluoranthene	14.95	252	94376m	4.58	ng	
89) Benzo(a)pyrene	15.27	252	298182	11.82	ng	97
90) Dibenzo(a,h)anthracene	16.67	278	51586	2.35	ng	# 88
91) Benzo(a,h,i)perylene	17.07	276	174444	7.81	ng	98

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

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