

Data Path : U:\HPCHEM1\BNA F\DATA\BF101816\
 Data File : BF090637.D
 Acq On : 18 Oct 2016 19:30
 Operator : UM/SJ
 Sample : H5289-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-2-2

Quant Time: Oct 19 01:59:52 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF101316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 17 19:19:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	238530	20.00	ng	-0.01
21) Naphthalene-d8	8.19	136	1057287	20.00	ng	-0.01
38) Acenaphthene-d10	9.95	164	521552	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	762871	20.00	ng	-0.01
75) Chrysene-d12	14.08	240	553331	20.00	ng	-0.01
86) Perylene-d12	15.54	264	462349	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	1547472	118.93	ng	0.00
7) Phenol-d6	6.54	99	2110127	109.14	ng	0.00
23) Nitrobenzene-d5	7.47	82	1364411	71.69	ng	-0.01
41) 2,4,6-Tribromophenol	10.74	330	380711	81.88	ng	-0.01
44) 2-Fluorobiphenyl	9.26	172	1822230	57.57	ng	-0.01
78) Terphenyl-d14	13.02	244	1452764	61.13	ng	-0.01
Target Compounds						Qvalue
48) Acenaphthylene	9.80	152	100893	2.14	ng	98
49) Dimethylphthalate	9.66	163	556724	16.45	ng	# 98
70) Phenanthrene	11.46	178	562079	13.80	ng	99
71) Anthracene	11.50	178	206046	4.70	ng	99
74) Fluoranthene	12.65	202	1185584	28.94	ng	93
77) Pyrene	12.88	202	1058731	28.89	ng	99
80) Benzo(a)anthracene	14.08	228	723147	22.96	ng	97
82) Chrysene	14.11	228	605026m	19.93	ng	
85) Indeno(1,2,3-cd)pyrene	16.99	276	242134	13.80	ng	# 91
87) Benzo(b)fluoranthene	15.12	252	937436m	30.98	ng	
88) Benzo(k)fluoranthene	15.14	252	216564m	6.65	ng	
89) Benzo(a)pyrene	15.48	252	452966	16.35	ng	96
90) Dibenzo(a,h)anthracene	17.00	278	60012	2.84	ng	94
91) Benzo(g,h,i)perylene	17.42	276	199711	9.92	ng	96

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

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