

Data Path : Z:\HPCHEM1\BNA F\DATA\BF112215\
 Data File : BF083179.D
 Acq On : 23 Nov 2015 1:05
 Operator : UM/IZ
 Sample : G4496-05
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 A19-COMP

Quant Time: Nov 23 06:07:37 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Nov 22 09:22:46 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	215895	20.00	ng	-0.01
21) Naphthalene-d8	7.94	136	787584	20.00	ng	-0.01
38) Acenaphthene-d10	9.68	164	383954	20.00	ng	-0.02
63) Phenanthrene-d10	11.16	188	662469	20.00	ng	-0.01
75) Chrysene-d12	13.78	240	500829	20.00	ng	-0.02
86) Perylene-d12	15.15	264	463128	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	1432613	100.32	ng	-0.02
7) Phenol-d6	6.32	99	2075990	112.36	ng	-0.02
23) Nitrobenzene-d5	7.22	82	1333615	77.37	ng	-0.01
41) 2,4,6-Tribromophenol	10.48	330	384262	97.82	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	1824116	66.25	ng	-0.01
78) Terphenyl-d14	12.74	244	1458484	68.58	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.42	163	519357	17.57	ng	99
70) Phenanthrene	11.19	178	129221	3.39	ng	98
74) Fluoranthene	12.36	202	294257	7.53	ng	98
77) Pyrene	12.59	202	232395	6.79	ng	98
80) Benzo(a)anthracene	13.77	228	114596	3.70	ng	99
82) Chrysene	13.81	228	95354	3.32	ng	97
85) Indeno(1,2,3-cd)pyrene	16.41	276	56292	2.15	ng	97
87) Benzo(b)fluoranthene	14.78	252	122082m	3.86	ng	
88) Benzo(k)fluoranthene	14.79	252	51374m	2.18	ng	
89) Benzo(a)pyrene	15.10	252	84825	3.25	ng	# 94
91) Benzo(g,h,i)perylene	16.79	276	52366	2.26	ng	# 94

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

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