

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120315\
 Data File : BF083463.D
 Acq On : 3 Dec 2015 16:40
 Operator : UM/IZ
 Sample : G4585-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP3-112415

Quant Time: Dec 03 17:29:21 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 03 17:02:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	174052	20.00	ng	-0.01
21) Naphthalene-d8	7.81	136	683114	20.00	ng	-0.01
38) Acenaphthene-d10	9.56	164	361694	20.00	ng	0.00
63) Phenanthrene-d10	11.09	188	637408	20.00	ng	0.00
75) Chrysene-d12	14.71	240	485390	20.00	ng	-0.01
86) Perylene-d12	16.77	264	384568	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	1600057	137.95	ng	0.01
7) Phenol-d6	6.20	99	1965685	123.81	ng	0.00
23) Nitrobenzene-d5	7.10	82	1419689	91.36	ng	0.00
41) 2,4,6-Tribromophenol	10.35	330	416655	114.75	ng	-0.01
44) 2-Fluorobiphenyl	8.90	172	2041806	80.45	ng	0.00
78) Terphenyl-d14	13.19	244	1888562	92.79	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.30	163	689544	23.46	ng	99
70) Phenanthrene	11.12	178	427581	11.34	ng	100
71) Anthracene	11.19	178	77623	2.04	ng	98
74) Fluoranthene	12.63	202	1064526	27.23	ng	96
77) Pyrene	12.93	202	1145048	33.78	ng	99
80) Benzo(a)anthracene	14.69	228	523068	17.59	ng	96
82) Chrysene	14.75	228	595874	20.74	ng	98
83) Bis(2-ethylhexyl)phthalate	14.82	149	47723	2.47	ng	97
85) Indeno(1,2,3-cd)pyrene	18.16	276	403095	19.26	ng	97
87) Benzo(b)fluoranthene	16.27	252	709278m	28.74	ng	
88) Benzo(k)fluoranthene	16.29	252	252407m	10.83	ng	
89) Benzo(a)pyrene	16.71	252	395737	18.67	ng	99
90) Dibenzo(a,h)anthracene	18.17	278	92203	4.99	ng	96
91) Benzo(a,h,i)perylene	18.52	276	431231	23.79	ng	# 95

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

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