

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120315\
 Data File : BF083472.D
 Acq On : 3 Dec 2015 21:20
 Operator : UM/IZ
 Sample : G4608-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CWC-MISC-3-20151130

Quant Time: Dec 04 01:22:20 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 01 01:17:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	174987	20.00	ng	-0.02
21) Naphthalene-d8	7.81	136	667164	20.00	ng	-0.02
38) Acenaphthene-d10	9.56	164	339467	20.00	ng	-0.02
63) Phenanthrene-d10	11.09	188	588292	20.00	ng	-0.03
75) Chrysene-d12	14.71	240	402337	20.00	ng	-0.06
86) Perylene-d12	16.77	264	356426	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.12	112	1087364	93.24	ng	-0.01
7) Phenol-d6	6.20	99	1520927	95.29	ng	-0.02
23) Nitrobenzene-d5	7.10	82	1116370	73.56	ng	-0.02
41) 2,4,6-Tribromophenol	10.35	330	270301	79.32	ng	-0.03
44) 2-Fluorobiphenyl	8.90	172	1524697	64.01	ng	-0.02
78) Terphenyl-d14	13.19	244	1138477	67.48	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.30	163	388597	14.09	ng	98
70) Phenanthrene	11.12	178	265073	7.61	ng	99
74) Fluoranthene	12.62	202	323549	8.97	ng	91
77) Pyrene	12.93	202	310216	11.04	ng	100
80) Benzo(a)anthracene	14.69	228	129686	5.26	ng	98
82) Chrysene	14.74	228	121474	5.10	ng	99
85) Indeno(1,2,3-cd)pyrene	18.15	276	56579	3.26	ng	97
87) Benzo(b)fluoranthene	16.26	252	133059m	5.82	ng	
88) Benzo(k)fluoranthene	16.29	252	44974m	2.08	ng	
89) Benzo(a)pyrene	16.69	252	98653	5.02	ng	# 91
91) Benzo(g,h,i)perylene	18.52	276	58249	3.47	ng	93

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

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