

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091115\
 Data File : BF081599.D
 Acq On : 12 Sep 2015 00:14
 Operator : UM/IZ
 Sample : G3526-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-13(0-0.16)

Quant Time: Sep 12 01:09:58 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 11 23:14:51 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.25	152	58022	20.00	ng	0.00
21) Naphthalene-d8	11.14	136	231825	20.00	ng	-0.01
38) Acenaphthene-d10	15.28	164	105047	20.00	ng	-0.01
63) Phenanthrene-d10	18.79	188	188208	20.00	ng	0.00
75) Chrysene-d12	23.40	240	116678	20.00	ng	-0.01
86) Perylene-d12	24.84	264	106652	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	339795	90.86	ng	0.02
7) Phenol-d6	7.67	99	500529	101.11	ng	-0.01
23) Nitrobenzene-d5	9.54	82	331494	68.99	ng	-0.01
41) 2,4,6-Tribromophenol	17.18	330	77739	83.11	ng	-0.02
44) 2-Fluorobiphenyl	13.80	172	518595	63.26	ng	-0.01
78) Terphenyl-d14	22.13	244	316626	63.17	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.85	163	95139	11.66	ng	# 97
70) Phenanthrene	18.84	178	34473	3.29	ng	94
74) Fluoranthene	21.44	202	58466	5.16	ng	89
77) Pyrene	21.80	202	53774	6.22	ng	98
80) Benzo(a)anthracene	23.38	228	30849	4.15	ng	# 90
82) Chrysene	23.42	228	23892	3.59	ng	93
85) Indeno(1,2,3-cd)pyrene	25.89	276	16751	3.43	ng	# 84
87) Benzo(b)fluoranthene	24.49	252	44228m	5.81	ng	
89) Benzo(a)pyrene	24.78	252	30257	4.69	ng	# 84
91) Benzo(g,h,i)perylene	26.18	276	15538	3.27	ng	# 78

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

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