

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060415\
 Data File : BF079722.D
 Acq On : 5 Jun 2015 2:47
 Operator : TP/IZ
 Sample : G2464-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-22AS

Quant Time: Jun 06 03:41:36 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 06 03:31:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	82842	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	328794	20.00	ng	0.00
38) Acenaphthene-d10	10.57	164	161968	20.00	ng	-0.01
63) Phenanthrene-d10	12.09	188	334057	20.00	ng	0.01
75) Chrysene-d12	15.02	240	373049	20.00	ng	0.03
86) Perylene-d12	16.96	264	376185	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	6.12	112	516938	107.00	ng	0.00
7) Phenol-d6	7.10	99	666720	106.01	ng	0.00
23) Nitrobenzene-d5	8.06	82	360088	66.68	ng	-0.01
41) 2,4,6-Tribromophenol	11.37	330	172691	118.73	ng	0.00
44) 2-Fluorobiphenyl	9.87	172	880112	78.67	ng	0.00
78) Terphenyl-d14	13.76	244	1060153	66.95	ng	0.02
Target Compounds						Qvalue
49) Dimethylphthalate	10.26	163	88889	7.12	ng	# 97
70) Phenanthrene	12.11	178	179484	9.24	ng	98
71) Anthracene	12.16	178	60910	3.17	ng	100
74) Fluoranthene	13.37	202	385370	18.28	ng	94
77) Pyrene	13.62	202	391568	16.89	ng	98
80) Benzo(a)anthracene	15.01	228	233732	10.25	ng	100
82) Chrysene	15.05	228	192902	9.04	ng	97
85) Indeno(1,2,3-cd)pyrene	18.96	276	152332	6.32	ng	93
87) Benzo(b)fluoranthene	16.38	252	253647m	10.45	ng	
88) Benzo(k)fluoranthene	16.40	252	116758m	5.04	ng	
89) Benzo(a)pyrene	16.88	252	207427	9.36	ng	97
90) Dibenzo(a,h)anthracene	18.98	278	50324	2.46	ng	# 92
91) Benzo(g,h,i)perylene	19.57	276	154549	7.38	ng	98

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

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