

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
 Data File : BF086409.D
 Acq On : 24 Apr 2016 00:14
 Operator : UM/SJ
 Sample : H2652-11 2X
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B1(0-7.5)-C

Manual Integrations
 APPROVED

sohil
 4/25/2016 12:36:15 PM

Quant Time: Apr 25 07:18:33 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 23 00:24:02 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	160236	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	629504	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	249463	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	431917	20.00	ng	0.00
75) Chrysene-d12	14.00	240	318770	20.00	ng	0.00
86) Perylene-d12	15.43	264	230103	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	696412	75.06	ng	0.01
7) Phenol-d6	6.48	99	915190	72.55	ng	0.00
23) Nitrobenzene-d5	7.40	82	522047	49.13	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	152208	67.40	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	923474	63.17	ng	0.00
78) Terphenyl-d14	12.94	244	668631	52.12	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	109995	6.10	ng	100
70) Phenanthrene	11.39	178	87691	4.04	ng	98
74) Fluoranthene	12.57	202	185884	7.94	ng	98
77) Pyrene	12.80	202	178039	8.01	ng	98
80) Benzo(a)anthracene	13.99	228	90874	5.49	ng	94
82) Chrysene	14.02	228	75523m	4.05	ng	
85) Indeno(1,2,3-cd)pyrene	16.81	276	37607	2.27	ng	96
87) Benzo(b)fluoranthene	15.01	252	74959m	5.93	ng	
88) Benzo(k)fluoranthene	15.03	252	41995m	3.39	ng	
89) Benzo(a)pyrene	15.36	252	63085	5.06	ng	# 78
91) Benzo(g,h,i)perylene	17.22	276	36633	3.45	ng	94

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

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