

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
 Data File : BF086408.D
 Acq On : 23 Apr 2016 23:45
 Operator : UM/SJ
 Sample : H2652-09
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B23(0-4.5)-C

Manual Integrations
 APPROVED

sohil
 4/25/2016 12:36:16 PM

Quant Time: Apr 25 07:16:37 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	172885	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	671317	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	276022	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	444209	20.00	ng	0.00
75) Chrysene-d12	14.00	240	348018	20.00	ng	0.00
86) Perylene-d12	15.41	264	250004	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1131233	113.00	ng	0.02
7) Phenol-d6	6.49	99	1456723	107.03	ng	0.01
23) Nitrobenzene-d5	7.41	82	859936	75.88	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	226452	90.63	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1298227	80.26	ng	0.00
78) Terphenyl-d14	12.95	244	926707	66.16	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.60	163	167776	8.41	ng	Qvalue 100
70) Phenanthrene	11.39	178	98219	4.40	ng	98
74) Fluoranthene	12.57	202	141442	5.88	ng	98
77) Pyrene	12.80	202	137425	5.66	ng	97
80) Benzo(a)anthracene	13.99	228	64846	3.59	ng	96
82) Chrysene	14.02	228	48553m	2.39	ng	
87) Benzo(b)fluoranthene	15.01	252	53805m	3.92	ng	
88) Benzo(k)fluoranthene	15.03	252	29769m	2.21	ng	
89) Benzo(a)pyrene	15.36	252	42593	3.15	ng	# 81
91) Benzo(g,h,i)perylene	17.21	276	24758	2.15	ng	# 92

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

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