

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042216\
 Data File : BF086392.D
 Acq On : 23 Apr 2016 16:02
 Operator : UM/SJ
 Sample : H2640-07
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WPG-B13(5-9)-C

Manual Integrations
 APPROVED

sohil
 4/25/2016 12:35:59 PM

Quant Time: Apr 25 05:44:17 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	180913	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	711181	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	325211	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	544869	20.00	ng	-0.01
75) Chrysene-d12	13.98	240	376742	20.00	ng	-0.01
86) Perylene-d12	15.40	264	313747	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1412365	134.82	ng	0.02
7) Phenol-d6	6.48	99	1710900	120.13	ng	0.00
23) Nitrobenzene-d5	7.40	82	1117263	93.06	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	357100	121.30	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	1799930	94.45	ng	0.00
78) Terphenyl-d14	12.94	244	1235236	81.46	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.60	163	237988	10.13	ng	Qvalue 100
70) Phenanthrene	11.38	178	245415	8.96	ng	99
71) Anthracene	11.44	178	76666	2.44	ng	100
74) Fluoranthene	12.57	202	266145	9.01	ng	95
77) Pyrene	12.80	202	233008	8.87	ng	99
80) Benzo(a)anthracene	13.98	228	87469	4.47	ng	99
82) Chrysene	14.02	228	80929m	3.67	ng	
87) Benzo(b)fluoranthene	15.00	252	62514m	3.63	ng	
88) Benzo(k)fluoranthene	15.01	252	43416m	2.57	ng	
89) Benzo(a)pyrene	15.35	252	60751	3.58	ng	# 88

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

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