

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051116\
 Data File : BF086916.D
 Acq On : 12 May 2016 4:49
 Operator : UM/SJ
 Sample : H2957-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP31-5FT

Manual Integrations
 APPROVED

sohil
 5/12/2016 9:09:11 AM

Quant Time: May 12 12:37:51 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 12 01:06:48 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	127950	20.00	ng	-0.01
21) Naphthalene-d8	7.92	136	487378	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	193941	20.00	ng	-0.01
63) Phenanthrene-d10	11.14	188	307601	20.00	ng	-0.01
75) Chrysene-d12	13.77	240	285780	20.00	ng	-0.01
86) Perylene-d12	15.14	264	224973	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	776343	102.76	ng	0.01
7) Phenol-d6	6.30	99	902033	89.33	ng	-0.01
23) Nitrobenzene-d5	7.21	82	614182	63.28	ng	-0.01
41) 2,4,6-Tribromophenol	10.46	330	162252	79.02	ng	-0.01
44) 2-Fluorobiphenyl	9.00	172	836753	72.51	ng	0.00
78) Terphenyl-d14	12.73	244	614953	45.66	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.40	163	116356	7.86	ng	# 98
70) Phenanthrene	11.16	178	39733	2.42	ng	98
74) Fluoranthene	12.35	202	123798	7.33	ng	93
77) Pyrene	12.58	202	113309	5.18	ng	98
80) Benzo(a)anthracene	13.76	228	69106	4.31	ng	# 90
82) Chrysene	13.79	228	53627m	3.29	ng	
87) Benzo(b)fluoranthene	14.77	252	65440m	4.48	ng	
88) Benzo(k)fluoranthene	14.78	252	29309m	2.45	ng	
89) Benzo(a)pyrene	15.09	252	46490	3.69	ng	# 91
91) Benzo(g,h,i)perylene	16.77	276	23803	2.20	ng	94

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

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