

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051116\
 Data File : BF086906.D
 Acq On : 12 May 2016 00:00
 Operator : UM/SJ
 Sample : H2952-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-1

Manual Integrations
 APPROVED

umangi
 5/12/2016 9:10:05 AM

Quant Time: May 12 01:42:27 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	149597	20.00	ng	-0.01
21) Naphthalene-d8	7.92	136	614231	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	284806	20.00	ng	-0.01
63) Phenanthrene-d10	11.14	188	486145	20.00	ng	-0.01
75) Chrysene-d12	13.77	240	319708	20.00	ng	-0.01
86) Perylene-d12	15.13	264	294537	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	950524	107.60	ng	0.00
7) Phenol-d6	6.30	99	1351460	114.48	ng	-0.01
23) Nitrobenzene-d5	7.20	82	896634	73.30	ng	-0.02
41) 2,4,6-Tribromophenol	10.46	330	307466	101.97	ng	-0.01
44) 2-Fluorobiphenyl	8.99	172	1267906	74.82	ng	-0.01
78) Terphenyl-d14	12.73	244	971567	64.48	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.39	163	233930	10.77	ng	99
70) Phenanthrene	11.17	178	56915	2.19	ng	99
74) Fluoranthene	12.35	202	146720	5.50	ng	90
77) Pyrene	12.57	202	133801	5.47	ng	98
80) Benzo(a)anthracene	13.76	228	81535	4.54	ng	98
82) Chrysene	13.79	228	55423	3.04	ng	98
87) Benzo(b)fluoranthene	14.77	252	92312m	4.82	ng	
89) Benzo(a)pyrene	15.07	252	59608	3.61	ng	# 93
91) Benzo(g,h,i)perylene	16.75	276	36908	2.61	ng	95

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

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