

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051116\
 Data File : BF086902.D
 Acq On : 11 May 2016 22:03
 Operator : UM/SJ
 Sample : H2894-03
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP24-7-8FT

Manual Integrations
 APPROVED

umangi
 5/12/2016 9:10:03 AM

Quant Time: May 12 01:28:47 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	154991	20.00	ng	-0.01
21) Naphthalene-d8	7.92	136	648319	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	303454	20.00	ng	-0.01
63) Phenanthrene-d10	11.14	188	498968	20.00	ng	-0.01
75) Chrysene-d12	13.77	240	376472	20.00	ng	-0.01
86) Perylene-d12	15.13	264	309798	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	1328298	145.14	ng	0.01
7) Phenol-d6	6.30	99	1617981	132.28	ng	-0.01
23) Nitrobenzene-d5	7.21	82	1218199	94.35	ng	-0.01
41) 2,4,6-Tribromophenol	10.46	330	409147	127.36	ng	-0.01
44) 2-Fluorobiphenyl	9.00	172	1593343	88.25	ng	0.00
78) Terphenyl-d14	12.73	244	1131016	63.74	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.40	163	222400	9.61	ng	99
74) Fluoranthene	12.35	202	139553	5.10	ng	95
77) Pyrene	12.58	202	130140	4.52	ng	98
80) Benzo(a)anthracene	13.76	228	73099	3.46	ng	93
82) Chrysene	13.79	228	59888m	2.79	ng	
87) Benzo(b)fluoranthene	14.77	252	72591m	3.61	ng	
89) Benzo(a)pyrene	15.09	252	50302	2.90	ng	# 83
91) Benzo(g,h,i)perylene	16.77	276	31687	2.13	ng	94

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

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