

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031116\
 Data File : BF085382.D
 Acq On : 11 Mar 2016 20:00
 Operator : UM/SJ
 Sample : H1757-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-STOCKPICE-COMP2

Manual Integrations
 APPROVED

UMANGI
 3/14/2016 10:45:24 AM

Quant Time: Mar 12 00:20:24 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 04 00:54:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	153312	20.00	ng	-0.05
21) Naphthalene-d8	8.22	136	630059	20.00	ng	-0.05
38) Acenaphthene-d10	9.98	164	287674	20.00	ng	-0.05
63) Phenanthrene-d10	11.46	188	516717	20.00	ng	-0.05
75) Chrysene-d12	14.11	240	300535	20.00	ng	-0.05
86) Perylene-d12	15.57	264	273811	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1457938	159.51	ng	-0.03
7) Phenol-d6	6.56	99	1756626	146.70	ng	-0.05
23) Nitrobenzene-d5	7.50	82	1037680	88.13	ng	-0.05
41) 2,4,6-Tribromophenol	10.77	330	368002	149.50	ng	-0.05
44) 2-Fluorobiphenyl	9.30	172	1899355	100.41	ng	-0.05
78) Terphenyl-d14	13.05	244	1363385	105.31	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.69	163	220534	9.99	ng	99
74) Fluoranthene	12.68	202	125445	4.62	ng	98
77) Pyrene	12.91	202	122519	5.42	ng	98
80) Benzo(a)anthracene	14.09	228	59743	3.32	ng	94
82) Chrysene	14.13	228	66417	4.00	ng	97
85) Indeno(1,2,3-cd)pyrene	17.02	276	42619	3.29	ng	# 91
87) Benzo(b)fluoranthene	15.15	252	105730m	5.79	ng	
88) Benzo(k)fluoranthene	15.17	252	31019m	2.11	ng	
89) Benzo(a)pyrene	15.50	252	44423	2.94	ng	# 90
91) Benzo(g,h,i)perylene	17.45	276	39200	3.02	ng	96

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

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