

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF031116\
 Data File : BF085381.D
 Acq On : 11 Mar 2016 19:32
 Operator : UM/SJ
 Sample : H1757-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TRACK-D-STOCKPICE-COMP1

Manual Integrations
 APPROVED

UMANGI
 3/14/2016 10:45:22 AM

Quant Time: Mar 14 04:55:28 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	142598	20.00	ng	-0.05
21) Naphthalene-d8	8.22	136	568922	20.00	ng	-0.05
38) Acenaphthene-d10	9.98	164	260740	20.00	ng	-0.05
63) Phenanthrene-d10	11.46	188	447333	20.00	ng	-0.05
75) Chrysene-d12	14.11	240	268996	20.00	ng	-0.05
86) Perylene-d12	15.57	264	262389	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1381272	162.48	ng	-0.03
7) Phenol-d6	6.56	99	1486667	133.48	ng	-0.05
23) Nitrobenzene-d5	7.50	82	975733	91.78	ng	-0.05
41) 2,4,6-Tribromophenol	10.77	330	322076	144.36	ng	-0.05
44) 2-Fluorobiphenyl	9.30	172	1757272	102.50	ng	-0.05
78) Terphenyl-d14	13.05	244	1235425	106.62	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.69	163	209956	10.49	ng	99
74) Fluoranthene	12.68	202	86483	3.68	ng	96
77) Pyrene	12.90	202	82880	4.09	ng	97
80) Benzo(a)anthracene	14.09	228	39076m	2.43	ng	
82) Chrysene	14.13	228	52193	3.51	ng	97
85) Indeno(1,2,3-cd)pyrene	17.02	276	38362	3.30	ng	95
87) Benzo(b)fluoranthene	15.13	252	85471m	4.89	ng	
89) Benzo(a)pyrene	15.50	252	35822	2.47	ng	# 90
91) Benzo(g,h,i)perylene	17.45	276	33895	2.72	ng	98

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

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