

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070616\
 Data File : BF088760.D
 Acq On : 7 Jul 2016 3:49
 Operator : UM/SJ
 Sample : H3879-11
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C2.1-04

Manual Integrations

umangi
 7/7/2016 6:10:33 PM

Quant Time: Jul 07 07:06:10 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 07 01:24:07 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	92847	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	381677	20.00	ng	0.00
38) Acenaphthene-d10	9.78	164	164186	20.00	ng	0.00
63) Phenanthrene-d10	11.26	188	262792	20.00	ng	0.00
75) Chrysene-d12	13.89	240	189404	20.00	ng	0.00
86) Perylene-d12	15.27	264	179599	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	566999	91.52	ng	0.00
7) Phenol-d6	6.39	99	667665	83.41	ng	0.00
23) Nitrobenzene-d5	7.31	82	425717	58.45	ng	-0.01
41) 2,4,6-Tribromophenol	10.57	330	152561	100.06	ng	0.00
44) 2-Fluorobiphenyl	9.11	172	766013	73.70	ng	-0.01
78) Terphenyl-d14	12.85	244	543907	63.96	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.51	163	110279	9.43	ng	99
70) Phenanthrene	11.28	178	53197	3.70	ng	99
71) Anthracene	11.32	178	39290	2.75	ng	97
74) Fluoranthene	12.47	202	225040	15.45	ng	92
77) Pyrene	12.68	202	218750	13.62	ng	98
80) Benzo(a)anthracene	13.87	228	107709m	8.83	ng	
82) Chrysene	13.91	228	128040	10.61	ng	97
85) Indeno(1,2,3-cd)pyrene	16.57	276	71653	7.72	ng	# 100
87) Benzo(b)fluoranthene	14.88	252	249001m	22.10	ng	
88) Benzo(k)fluoranthene	14.90	252	25364m	2.59	ng	
89) Benzo(a)pyrene	15.21	252	85066	8.92	ng	96
90) Dibenzo(a,h)anthracene	16.58	278	17254	2.17	ng	# 87
91) Benzo(a,h,i)perylene	16.97	276	65778	7.98	ng	98

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

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