

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071316\
 Data File : BF088960.D
 Acq On : 13 Jul 2016 16:16
 Operator : UM/SJ
 Sample : H3876-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 5-A-LOCATION-5-0-5FT

Manual Integrations

APPROVED
 Sohil
 7/14/2016 4:19:45 AM

Quant Time: Jul 14 01:25:16 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 14 01:23:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	71540m	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	263696	20.00	ng	-0.01
38) Acenaphthene-d10	9.72	164	119426	20.00	ng	-0.01
63) Phenanthrene-d10	11.20	188	195115	20.00	ng	-0.01
75) Chrysene-d12	13.83	240	126161	20.00	ng	-0.01
86) Perylene-d12	15.20	264	132430	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	393214	82.38	ng	0.00
7) Phenol-d6	6.34	99	495454	80.33	ng	-0.01
23) Nitrobenzene-d5	7.26	82	332176	66.01	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	110662	99.79	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	596343	78.87	ng	0.00
78) Terphenyl-d14	12.79	244	348069	61.45	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.45	163	88523	10.40	ng	100
70) Phenanthrene	11.22	178	27219	2.55	ng	99
74) Fluoranthene	12.41	202	53212	4.92	ng	95
77) Pyrene	12.64	202	43720	4.09	ng	98
80) Benzo(a)anthracene	13.82	228	22776	2.80	ng	92
87) Benzo(b)fluoranthene	14.82	252	23094m	2.78	ng	
89) Benzo(a)pyrene	15.14	252	16625	2.36	ng	96

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

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