

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
 Data File : BF088585.D
 Acq On : 1 Jul 2016 20:13
 Operator : UM/SJ
 Sample : H3836-08 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C1.1-01

Manual Integrations
 APPROVED

sohil
 7/5/2016 7:31:14 PM

Quant Time: Jul 02 03:25:28 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 30 18:01:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	132433	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	501264	20.00	ng	-0.01
38) Acenaphthene-d10	9.84	164	209975	20.00	ng	-0.01
63) Phenanthrene-d10	11.31	188	328977	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	262997	20.00	ng	-0.01
86) Perylene-d12	15.35	264	220545	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	389772	44.11	ng	-0.01
7) Phenol-d6	6.43	99	533108	46.69	ng	-0.01
23) Nitrobenzene-d5	7.37	82	379238	39.65	ng	-0.01
41) 2,4,6-Tribromophenol	10.63	330	70391	36.10	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	556858	41.89	ng	-0.01
78) Terphenyl-d14	12.90	244	400015	33.88	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.56	163	48360	3.23	ng	99
70) Phenanthrene	11.34	178	66139	3.68	ng	99
74) Fluoranthene	12.53	202	97654	5.36	ng	97
77) Pyrene	12.75	202	76849	3.45	ng	98
80) Benzo(a)anthracene	13.93	228	43067	2.54	ng	95
87) Benzo(b)fluoranthene	14.95	252	45854m	3.31	ng	
89) Benzo(a)pyrene	15.28	252	25439	2.17	ng	# 84
91) Benzo(g,h,i)perylene	17.09	276	22150	2.19	ng	97

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

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