

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070116\
 Data File : BF088570.D
 Acq On : 1 Jul 2016 12:54
 Operator : UM/SJ
 Sample : H3808-06 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-COMP

Manual Integrations
 APPROVED

sohil
 7/5/2016 7:30:51 PM

Quant Time: Jul 01 14:20:25 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	143373	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	550015	20.00	ng	-0.01
38) Acenaphthene-d10	9.84	164	239514	20.00	ng	-0.01
63) Phenanthrene-d10	11.33	188	459559	20.00	ng	0.00
75) Chrysene-d12	13.94	240	306894	20.00	ng	-0.01
86) Perylene-d12	15.35	264	311443	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	435551	45.53	ng	-0.01
7) Phenol-d6	6.43	99	596724	48.27	ng	-0.01
23) Nitrobenzene-d5	7.37	82	400292	38.14	ng	-0.01
41) 2,4,6-Tribromophenol	10.63	330	118971	53.49	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	563832	37.18	ng	-0.01
78) Terphenyl-d14	12.90	244	451576	32.77	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.56	163	57968	3.40	ng	Qvalue 98
70) Phenanthrene	11.35	178	84253	3.35	ng	97
74) Fluoranthene	12.53	202	193182	7.59	ng	98
77) Pyrene	12.75	202	162161	6.23	ng	99
80) Benzo(a)anthracene	13.93	228	87699	4.44	ng	99
82) Chrysene	13.97	228	64303m	3.29	ng	
85) Indeno(1,2,3-cd)pyrene	16.69	276	37238	2.48	ng	# 100
87) Benzo(b)fluoranthene	14.95	252	85665m	4.38	ng	
88) Benzo(k)fluoranthene	14.96	252	37488m	2.20	ng	
89) Benzo(a)pyrene	15.29	252	63057	3.81	ng	96
91) Benzo(g,h,i)perylene	17.09	276	37240	2.61	ng	95

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

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