

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062016\
 Data File : BF088209.D
 Acq On : 21 Jun 2016 9:57
 Operator : UM/SJ
 Sample : H3474-03
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S7-B4(6-12)C

Manual Integrations
 APPROVED

sohil
 6/21/2016 4:22:28 PM

Quant Time: Jun 21 13:01:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 20 17:39:32 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	102943	20.00	ng	-0.01
21) Naphthalene-d8	7.95	136	397889	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	170089	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	262119	20.00	ng	0.00
75) Chrysene-d12	13.81	240	201712	20.00	ng	0.00
86) Perylene-d12	15.19	264	229160	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	693990	115.25	ng	0.01
7) Phenol-d6	6.33	99	859581	117.79	ng	0.00
23) Nitrobenzene-d5	7.23	82	478568	70.23	ng	-0.01
41) 2,4,6-Tribromophenol	10.49	330	151246	84.21	ng	0.00
44) 2-Fluorobiphenyl	9.03	172	705024	63.73	ng	-0.01
78) Terphenyl-d14	12.75	244	474142	53.49	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.43	163	171977	13.96	ng	99
70) Phenanthrene	11.20	178	129177	9.39	ng	98
71) Anthracene	11.26	178	59207	4.27	ng	99
74) Fluoranthene	12.39	202	432328	29.78	ng	98
77) Pyrene	12.62	202	380764	25.44	ng	98
80) Benzo(a)anthracene	13.79	228	306182	26.64	ng	99
82) Chrysene	13.83	228	212892	19.69	ng	98
85) Indeno(1,2,3-cd)pyrene	16.47	276	140927	14.03	ng	# 100
87) Benzo(b)fluoranthene	14.81	252	361522m	22.63	ng	
88) Benzo(k)fluoranthene	14.82	252	97953m	8.13	ng	
89) Benzo(a)pyrene	15.13	252	252600	19.55	ng	97
90) Dibenzo(a,h)anthracene	16.48	278	47877	3.99	ng	# 86
91) Benzo(a,h,i)perylene	16.86	276	121381	9.83	ng	97

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

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