

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062317\
 Data File : BF096169.D
 Acq On : 23 Jun 2017 20:25
 Operator : SJ/MA
 Sample : I3784-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P.56-B-COMP

Manual Integrations
 APPROVED

mohammad
 6/28/2017 1:11:24 PM

Quant Time: Jun 24 00:43:32 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 20 16:05:47 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.26	152	70515	20.00	ng	-0.04
21) Naphthalene-d8	9.29	136	295107	20.00	ng	-0.04
38) Acenaphthene-d10	12.11	164	133976	20.00	ng	-0.04
63) Phenanthrene-d10	14.50	188	231089	20.00	ng	-0.04
75) Chrysene-d12	18.24	240	153174	20.00	ng	-0.03
86) Perylene-d12	19.92	264	138023	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.15	112	558342	126.17	ng	-0.04
7) Phenol-d6	6.84	99	675451	127.50	ng	-0.04
23) Nitrobenzene-d5	8.19	82	393938	82.90	ng	-0.04
41) 2,4,6-Tribromophenol	13.42	330	141202	119.12	ng	-0.04
44) 2-Fluorobiphenyl	11.07	172	619701	64.37	ng	-0.04
78) Terphenyl-d14	16.90	244	439725	71.24	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	11.77	163	144753	14.23	ng	98
70) Phenanthrene	14.54	178	77178	5.79	ng	96
74) Fluoranthene	16.34	202	102130	7.74	ng	97
77) Pyrene	16.64	202	94880	8.30	ng	# 95
80) Benzo(a)anthracene	18.24	228	34559	3.88	ng	96
82) Chrysene	18.29	228	35913	4.04	ng	96
85) Indeno(1,2,3-cd)pyrene	21.10	276	19221	2.52	ng	100
87) Benzo(b)fluoranthene	19.53	252	35347m	4.09	ng	
89) Benzo(a)pyrene	19.87	252	29275	3.69	ng	# 91
91) Benzo(g,h,i)perylene	21.43	276	17285	2.52	ng	# 88

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

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