

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061017\
 Data File : BF095866.D
 Acq On : 11 Jun 2017 1:01
 Operator : SJ/MA
 Sample : I3594-07 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-060917-C

Manual Integrations
 APPROVED

mohammad
 6/12/2017 1:28:43 PM

Quant Time: Jun 12 06:43:50 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:15:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	59703	20.00	ng	-0.03
21) Naphthalene-d8	9.40	136	235456	20.00	ng	-0.04
38) Acenaphthene-d10	12.22	164	95937	20.00	ng	-0.04
63) Phenanthrene-d10	14.62	188	151065	20.00	ng	-0.03
75) Chrysene-d12	18.34	240	137618	20.00	ng	-0.02
86) Perylene-d12	20.01	264	95177	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	76759	20.49	ng	0.00
7) Phenol-d6	6.96	99	96904	21.60	ng	0.00
23) Nitrobenzene-d5	8.29	82	53339	14.07	ng	-0.03
41) 2,4,6-Tribromophenol	13.53	330	15418	18.16	ng	-0.02
44) 2-Fluorobiphenyl	11.18	172	121942	17.69	ng	-0.04
78) Terphenyl-d14	16.99	244	85934	15.50	ng	-0.02
Target Compounds						
49) Dimethylphthalate	11.88	163	14998	2.06	ng	Qvalue # 95
70) Phenanthrene	14.66	178	83446	9.57	ng	96
71) Anthracene	14.74	178	27425	3.01	ng	94
74) Fluoranthene	16.43	202	112755	13.07	ng	89
77) Pyrene	16.74	202	122622	11.94	ng	# 95
80) Benzo(a)anthracene	18.33	228	58510m	7.31	ng	
82) Chrysene	18.37	228	53822	6.73	ng	98
85) Indeno(1,2,3-cd)pyrene	21.20	276	25104	3.67	ng	# 91
87) Benzo(b)fluoranthene	19.61	252	63498m	10.65	ng	
88) Benzo(k)fluoranthene	19.63	252	17921m	3.15	ng	
89) Benzo(a)pyrene	19.96	252	43399	7.92	ng	# 90
91) Benzo(a,h,i)perylene	21.53	276	23144	4.89	ng	91

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

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