

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051917\
 Data File : BF095279.D
 Acq On : 19 May 2017 23:47
 Operator : SJ/MA
 Sample : I3202-06
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-6COMP

Manual Integrations
 APPROVED

mohammad
 5/22/2017 1:29:57 PM

Quant Time: May 20 02:15:14 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 17:07:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	94571	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	343663	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	191492	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	344225	20.00	ng	0.00
75) Chrysene-d12	18.69	240	226517	20.00	ng	0.00
86) Perylene-d12	20.36	264	164843	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.73	112	581080	102.44	ng	0.00
7) Phenol-d6	7.30	99	709848	104.30	ng	-0.01
23) Nitrobenzene-d5	8.66	82	432531	79.36	ng	-0.01
41) 2,4,6-Tribromophenol	13.91	330	174587	111.33	ng	0.00
44) 2-Fluorobiphenyl	11.55	172	824615	61.98	ng	0.00
78) Terphenyl-d14	17.34	244	734978	71.47	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	12.25	163	139975	8.90	ng	99
70) Phenanthrene	15.05	178	145842	7.37	ng	98
74) Fluoranthene	16.80	202	196591	9.69	ng	97
77) Pyrene	17.10	202	199156	11.76	ng	96
80) Benzo(a)anthracene	18.68	228	73572	5.58	ng	99
82) Chrysene	18.72	228	79002m	6.20	ng	
85) Indeno(1,2,3-cd)pyrene	21.72	276	39919	3.86	ng	# 89
87) Benzo(b)fluoranthene	19.96	252	72122m	7.08	ng	
88) Benzo(k)fluoranthene	19.99	252	30148m	3.07	ng	
89) Benzo(a)pyrene	20.31	252	65288	6.95	ng	99
91) Benzo(g,h,i)perylene	22.11	276	40322	5.29	ng	95

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

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