

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051917\
 Data File : BF095304.D
 Acq On : 20 May 2017 14:42
 Operator : SJ/MA
 Sample : I3186-01 10X
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-S3-WSW

Manual Integrations
 APPROVED

mohammad
 5/22/2017 1:30:42 PM

Quant Time: May 22 04:39:40 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	79073	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	330946	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	142402	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	198197	20.00	ng	0.00
75) Chrysene-d12	18.69	240	184437	20.00	ng	0.00
86) Perylene-d12	20.37	264	143154	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.83	112	21928	4.62	ng	0.09
7) Phenol-d6	7.41	99	44365	7.80	ng	0.09
23) Nitrobenzene-d5	8.71	82	24341	4.64	ng	0.04
41) 2,4,6-Tribromophenol	13.94	330	7819	6.70	ng	0.02
44) 2-Fluorobiphenyl	11.57	172	74925	7.57	ng	0.00
78) Terphenyl-d14	17.34	244	43761	5.23	ng	-0.01
Target Compounds						
70) Phenanthrene	15.05	178	111865	9.82	ng	98
71) Anthracene	15.15	178	24169	2.08	ng	97
74) Fluoranthene	16.79	202	182644	15.64	ng	99
77) Pyrene	17.10	202	156997	11.38	ng	97
80) Benzo(a)anthracene	18.68	228	77079	7.18	ng	96
82) Chrysene	18.72	228	79976	7.71	ng	97
85) Indeno(1,2,3-cd)pyrene	21.71	276	25580	3.03	ng	98
87) Benzo(b)fluoranthene	19.96	252	85684m	9.69	ng	
88) Benzo(k)fluoranthene	19.98	252	31591m	3.70	ng	
89) Benzo(a)pyrene	20.31	252	55136	6.75	ng	99
91) Benzo(g,h,i)perylene	22.11	276	22451m	3.39	ng	

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

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