

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051817\
 Data File : BF095263.D
 Acq On : 19 May 2017 13:42
 Operator : SJ/MA
 Sample : I3186-09
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-S4-SSW

Manual Integrations
 APPROVED

mohammad
 5/19/2017 3:41:03 PM

Quant Time: May 19 15:15:28 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	94676	20.00	ng	0.00
21) Naphthalene-d8	9.79	136	420142	20.00	ng	0.00
38) Acenaphthene-d10	12.62	164	181273	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	279151	20.00	ng	0.00
75) Chrysene-d12	18.69	240	183529	20.00	ng	0.00
86) Perylene-d12	20.36	264	184397	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.74	112	555371	97.80	ng	0.00
7) Phenol-d6	7.31	99	648066	95.11	ng	0.00
23) Nitrobenzene-d5	8.67	82	399028	59.89	ng	0.00
41) 2,4,6-Tribromophenol	13.92	330	128979	86.88	ng	0.00
44) 2-Fluorobiphenyl	11.55	172	775701	61.59	ng	0.00
78) Terphenyl-d14	17.34	244	471658	56.60	ng	-0.01
Target Compounds						
49) Dimethylphthalate	12.26	163	180202	12.11	ng	98
70) Phenanthrene	15.06	178	82503	5.14	ng	97
74) Fluoranthene	16.80	202	159437	9.69	ng	99
77) Pyrene	17.11	202	149041	10.86	ng	96
80) Benzo(a)anthracene	18.68	228	59734	5.59	ng	98
82) Chrysene	18.73	228	63586	6.16	ng	98
85) Indeno(1,2,3-cd)pyrene	21.71	276	37495	4.47	ng	96
87) Benzo(b)fluoranthene	19.97	252	77076m	6.76	ng	
88) Benzo(k)fluoranthene	19.99	252	32023m	2.91	ng	
89) Benzo(a)pyrene	20.31	252	64082	6.09	ng	96
91) Benzo(g,h,i)perylene	22.10	276	40173m	4.71	ng	

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

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