

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051617\
 Data File : BF095182.D
 Acq On : 17 May 2017 5:38
 Operator : SJ/MA
 Sample : I3153-07
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DS-4-4

Quant Time: May 17 08:57:13 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 03 17:24:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.78	152	90669	20.00	ng	0.05
21) Naphthalene-d8	9.79	136	349581	20.00	ng	0.04
38) Acenaphthene-d10	12.62	164	116131	20.00	ng	0.04
63) Phenanthrene-d10	15.02	188	176730	20.00	ng	0.04
75) Chrysene-d12	18.69	240	175498	20.00	ng	0.04
86) Perylene-d12	20.36	264	131547	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.75	112	616851	113.43	ng	0.07
7) Phenol-d6	7.32	99	661078	101.31	ng	0.04
23) Nitrobenzene-d5	8.67	82	415071	74.87	ng	0.04
41) 2,4,6-Tribromophenol	13.92	330	92453	97.21	ng	0.04
44) 2-Fluorobiphenyl	11.56	172	615421	76.28	ng	0.03
78) Terphenyl-d14	17.34	244	402815	50.55	ng	0.03
Target Compounds						Qvalue
49) Dimethylphthalate	12.25	163	81988	8.60	ng	# 98
70) Phenanthrene	15.07	178	57650	5.68	ng	98
74) Fluoranthene	16.81	202	117241	11.26	ng	97
77) Pyrene	17.11	202	113354	8.64	ng	98
80) Benzo(a)anthracene	18.68	228	53323m	5.22	ng	
82) Chrysene	18.72	228	67517	6.84	ng	96
83) Bis(2-ethylhexyl)phthalate	18.74	149	114113	13.12	ng	99
85) Indeno(1,2,3-cd)pyrene	21.71	276	28809	3.59	ng	94
87) Benzo(b)fluoranthene	19.96	252	68690m	8.45	ng	
88) Benzo(k)fluoranthene	19.99	252	29172m	3.72	ng	
89) Benzo(a)pyrene	20.32	252	41901	5.59	ng	# 95
91) Benzo(a,h,i)perylene	22.10	276	31786	5.23	ng	96

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

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