

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052717\
 Data File : BF095527.D
 Acq On : 28 May 2017 9:19
 Operator : SJ/MA
 Sample : I3339-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Q3-WSW

Manual Integrations
 APPROVED

mohammad
 5/31/2017 1:05:15 PM

Quant Time: May 29 04:34:54 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 26 16:50:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	91649	20.00	ng	0.00
21) Naphthalene-d8	9.77	136	350122	20.00	ng	0.00
38) Acenaphthene-d10	12.60	164	141622	20.00	ng	0.00
63) Phenanthrene-d10	15.00	188	202803	20.00	ng	0.00
75) Chrysene-d12	18.67	240	199133	20.00	ng	0.00
86) Perylene-d12	20.35	264	144063	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	599324	109.02	ng	0.00
7) Phenol-d6	7.28	99	711656	107.90	ng	-0.01
23) Nitrobenzene-d5	8.64	82	396927	71.49	ng	-0.01
41) 2,4,6-Tribromophenol	13.90	330	104659	90.24	ng	0.00
44) 2-Fluorobiphenyl	11.53	172	735396	74.74	ng	0.00
78) Terphenyl-d14	17.31	244	420047	46.46	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	12.24	163	268988	23.14	ng	99
70) Phenanthrene	15.04	178	101220	8.69	ng	95
74) Fluoranthene	16.78	202	196531	16.44	ng	98
77) Pyrene	17.08	202	192905	12.95	ng	97
80) Benzo(a)anthracene	18.66	228	100693m	8.69	ng	
82) Chrysene	18.71	228	107183	9.56	ng	98
85) Indeno(1,2,3-cd)pyrene	21.71	276	34635	3.81	ng	94
87) Benzo(b)fluoranthene	19.95	252	99583m	11.19	ng	
88) Benzo(k)fluoranthene	19.98	252	36808m	4.29	ng	
89) Benzo(a)pyrene	20.30	252	69597	8.47	ng	# 93
91) Benzo(g,h,i)perylene	22.09	276	37906	5.69	ng	100

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

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