

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052617\
 Data File : BF095459.D
 Acq On : 26 May 2017 14:14
 Operator : SJ/MA
 Sample : I3240-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP2-WC8(0-8)

Manual Integrations
 APPROVED

mohammad
 5/30/2017 1:35:56 PM

Quant Time: May 26 17:07:27 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	95378	20.00	ng	0.00
21) Naphthalene-d8	9.76	136	396040	20.00	ng	0.00
38) Acenaphthene-d10	12.59	164	170483	20.00	ng	0.00
63) Phenanthrene-d10	15.00	188	256779	20.00	ng	0.00
75) Chrysene-d12	18.68	240	188471	20.00	ng	0.00
86) Perylene-d12	20.36	264	173819	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	732009	127.96	ng	0.00
7) Phenol-d6	7.29	99	864712	125.98	ng	0.00
23) Nitrobenzene-d5	8.64	82	499167	79.48	ng	-0.01
41) 2,4,6-Tribromophenol	13.90	330	143677	102.91	ng	0.00
44) 2-Fluorobiphenyl	11.53	172	885518	74.76	ng	0.00
78) Terphenyl-d14	17.32	244	527334	61.63	ng	0.00
Target Compounds						Qvalue
10) Phenol	7.31	94	47640	6.59	ng	86
20) 3+4-Methylphenols	8.57	107	48312	6.67	ng	# 63
27) 2,4-Dimethylphenol	9.39	122	12850	2.24	ng	98
31) Naphthalene	9.81	128	3293826	165.48	ng	99
37) 2-Methylnaphthalene	10.92	142	976969	74.79	ng	98
45) 1,1'-Biphenyl	11.68	154	329350	22.95	ng	98
48) Acenaphthylene	12.37	152	39096	2.15	ng	# 82
49) Dimethylphthalate	12.22	163	161980	11.57	ng	98
51) Acenaphthene	12.65	154	623420	57.50	ng	99
54) Dibenzofuran	12.94	168	1169646	77.95	ng	99
57) Fluorene	13.49	166	420286	32.21	ng	98
70) Phenanthrene	15.07	178	4257457	288.57	ng	100
71) Anthracene	15.12	178	227709m	15.11	ng	
72) Carbazole	15.41	167	355512	25.93	ng	98
74) Fluoranthene	16.81	202	2930143	193.61	ng	99
77) Pyrene	17.11	202	2590531	183.78	ng	100
80) Benzo(a)anthracene	18.67	228	912635	83.20	ng	99
82) Chrysene	18.72	228	1239871	116.90	ng	97
85) Indeno(1,2,3-cd)pyrene	21.69	276	433475	50.33	ng	93
87) Benzo(b)fluoranthene	19.96	252	1363320m	126.93	ng	
88) Benzo(k)fluoranthene	19.98	252	377073m	36.40	ng	
89) Benzo(a)pyrene	20.31	252	834514	84.20	ng	98
90) Dibenzo(a,h)anthracene	21.68	278	105827m	12.93	ng	
91) Benzo(g,h,i)perylene	22.07	276	410912	51.16	ng	98

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052617\
Data File : BF095459.D
Acq On : 26 May 2017 14:14
Operator : SJ/MA
Sample : I3240-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
EP2-WC8(0-8)

Manual Integrations
APPROVED

mohammad
5/30/2017 1:35:56 PM

Quant Time: May 26 17:07:27 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

