

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020217\
 Data File : BF092741.D
 Acq On : 2 Feb 2017 18:04
 Operator : SJ/MA
 Sample : I1659-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CENTER-COMP

Manual Integrations
 APPROVED

mohammad
 2/3/2017 1:22:25 PM

Quant Time: Feb 03 00:29:45 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 30 14:50:05 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	166525	20.00	ng	-0.03
21) Naphthalene-d8	8.22	136	600201	20.00	ng	-0.03
38) Acenaphthene-d10	9.97	164	297377	20.00	ng	-0.03
63) Phenanthrene-d10	11.46	188	484409	20.00	ng	-0.03
75) Chrysene-d12	14.10	240	331207	20.00	ng	-0.05
86) Perylene-d12	15.56	264	303803	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	830087	85.52	ng	-0.03
7) Phenol-d6	6.57	99	1081135	86.57	ng	-0.03
23) Nitrobenzene-d5	7.50	82	557965	51.77	ng	-0.03
41) 2,4,6-Tribromophenol	10.76	330	176822	82.21	ng	-0.05
44) 2-Fluorobiphenyl	9.30	172	947062	57.70	ng	-0.03
78) Terphenyl-d14	13.05	244	792669	56.23	ng	-0.03
Target Compounds						
10) Phenol	6.59	94	46739	3.20	ng	Qvalue # 65
49) Dimethylphthalate	9.69	163	119421	5.96	ng	98
51) Acenaphthene	10.01	154	43195	2.48	ng	96
57) Fluorene	10.52	166	47518	2.65	ng	99
70) Phenanthrene	11.48	178	608128	21.91	ng	97
71) Anthracene	11.54	178	146232	5.25	ng	95
72) Carbazole	11.69	167	80624	3.03	ng	98
74) Fluoranthene	12.67	202	773113	27.01	ng	98
77) Pyrene	12.90	202	674103	27.20	ng	99
80) Benzo(a)anthracene	14.09	228	313656m	15.48	ng	
82) Chrysene	14.12	228	287345m	15.15	ng	
83) Bis(2-ethylhexyl)phthalate	14.08	149	302831	22.00	ng	# 95
85) Indeno(1,2,3-cd)pyrene	17.00	276	168875	11.50	ng	99
87) Benzo(b)fluoranthene	15.13	252	363443m	18.18	ng	
88) Benzo(k)fluoranthene	15.15	252	82918m	4.80	ng	
89) Benzo(a)pyrene	15.49	252	237867	13.75	ng	99
90) Dibenzo(a,h)anthracene	17.01	278	38692m	2.77	ng	
91) Benzo(g,h,i)perylene	17.45	276	157459	11.15	ng	# 93

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

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