

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091317\
 Data File : BF098485.D
 Acq On : 13 Sep 2017 13:01
 Operator : SJ/JU
 Sample : I5140-19
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-13-16-32

Manual Integrations
 APPROVED

Sohil
 9/14/2017 5:52:17 PM

Quant Time: Sep 13 13:46:46 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	180443	20.00	ng	-0.02
21) Naphthalene-d8	7.93	136	743816	20.00	ng	-0.02
38) Acenaphthene-d10	9.69	164	327489	20.00	ng	-0.02
63) Phenanthrene-d10	11.17	188	627168	20.00	ng	-0.02
75) Chrysene-d12	13.80	240	421250	20.00	ng	-0.02
86) Perylene-d12	15.20	264	338167	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1402121	114.89	ng	0.00
7) Phenol-d6	6.30	99	1708310	110.25	ng	-0.02
23) Nitrobenzene-d5	7.22	82	979737	73.43	ng	-0.02
41) 2,4,6-Tribromophenol	10.47	330	284777	130.29	ng	-0.02
44) 2-Fluorobiphenyl	9.01	172	1291112	66.82	ng	-0.02
78) Terphenyl-d14	12.75	244	982966	50.37	ng	-0.02
Target Compounds						
10) Phenol	6.32	94	62112	3.451	ng	97
20) 3+4-Methylphenols	7.08	107	91259	6.608	ng	# 57
31) Naphthalene	7.96	128	157817	4.292	ng	99
49) Dimethylphthalate	9.41	163	199287	7.861	ng	98
70) Phenanthrene	11.19	178	316003	9.504	ng	98
74) Fluoranthene	12.37	202	353095	10.609	ng	99
77) Pyrene	12.60	202	307916	9.266	ng	99
80) Benzo(a)anthracene	13.79	228	114528	4.637	ng	96
82) Chrysene	13.82	228	116207	4.655	ng	97
85) Indeno(1,2,3-cd)pyrene	16.53	276	45266	2.584	ng	95
87) Benzo(b)fluoranthene	14.80	252	109597m	5.154	ng	
89) Benzo(a)pyrene	15.14	252	78031	4.119	ng	# 96
91) Benzo(a,h,i)perylene	16.93	276	48395	3.492	ng	99

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

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