

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091217\
 Data File : BF098466.D
 Acq On : 12 Sep 2017 22:38
 Operator : SJ/JU
 Sample : I5140-09
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RL-10-0-10.5

Manual Integrations
 APPROVED

Sohil
 9/13/2017 5:00:18 PM

Quant Time: Sep 13 02:19:47 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	6.66	152	201649	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	782593	20.00	ng	0.00
38) Acenaphthene-d10	9.69	164	301743	20.00	ng	-0.01
63) Phenanthrene-d10	11.17	188	464679	20.00	ng	-0.01
75) Chrysene-d12	13.81	240	370644	20.00	ng	0.00
86) Perylene-d12	15.21	264	287121	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	1306330	95.78	ng	0.00
7) Phenol-d6	6.32	99	1590336	91.84	ng	0.00
23) Nitrobenzene-d5	7.23	82	966065	68.82	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	188000	93.35	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1207736	67.84	ng	-0.01
78) Terphenyl-d14	12.76	244	771532	44.93	ng	-0.01
Target Compounds						
10) Phenol	6.33	94	42177	2.097	ng	97
49) Dimethylphthalate	9.42	163	199875	8.557	ng	99
70) Phenanthrene	11.20	178	125471	5.093	ng	98
74) Fluoranthene	12.39	202	261933	10.622	ng	98
77) Pyrene	12.62	202	230079	7.869	ng	99
80) Benzo(a)anthracene	13.80	228	112885	5.195	ng	99
82) Chrysene	13.83	228	106942	4.869	ng	94
85) Indeno(1,2,3-cd)pyrene	16.55	276	40231	2.611	ng	97
87) Benzo(b)fluoranthene	14.82	252	120233m	6.660	ng	
88) Benzo(k)fluoranthene	14.84	252	41481m	2.461	ng	
89) Benzo(a)pyrene	15.15	252	80112	4.981	ng	# 89
91) Benzo(a,h,i)perylene	16.96	276	39970	3.397	ng	95

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

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