

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
 Data File : BF111295.D
 Acq On : 12 Dec 2018 00:12
 Operator : JU/SJ
 Sample : J6148-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-MH-J

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:30:12 PM

Quant Time: Dec 12 01:17:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 01:15:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	267201	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	1053101	20.00	ng	-0.01
39) Acenaphthene-d10	9.83	164	473857	20.00	ng	-0.01
64) Phenanthrene-d10	11.32	188	760300	20.00	ng	0.00
76) Chrysene-d12	13.96	240	522667	20.00	ng	0.00
87) Perylene-d12	15.40	264	531844	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1557202	92.03	ng	0.00
7) Phenol-d6	6.43	99	1981491	94.23	ng	0.00
23) Nitrobenzene-d5	7.35	82	1253123	66.74	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	371547	88.52	ng	-0.01
45) 2-Fluorobiphenyl	9.16	172	1890127	64.30	ng	0.00
79) Terphenyl-d14	12.90	244	1354288	62.15	ng	0.00
Target Compounds						
10) Phenol	6.44	94	167872	6.827	ng	90
31) Naphthalene	8.10	128	109264	2.377	ng	96
37) 2-Methylnaphthalene	8.79	142	62346m	2.043	ng	
38) 1-Methylnaphthalene	8.89	142	64976	2.174	ng	# 98
49) Acenaphthylene	9.69	152	610093	13.554	ng	94
50) Dimethylphthalate	9.55	163	159400	4.559	ng	99
58) Fluorene	10.37	166	120182	4.041	ng	92
71) Phenanthrene	11.34	178	1206450	31.460	ng	95
72) Anthracene	11.39	178	351065	8.750	ng	94
73) Carbazole	11.54	167	150778	3.594	ng	96
75) Fluoranthene	12.53	202	1969703	56.227	ng	99
78) Pvrene	12.76	202	2351410	62.956	ng	98
80) Butylbenzylphthalate	13.39	149	2051494	105.158	ng	99
81) Benzo(a)anthracene	13.94	228	1240684m	42.549	ng	
83) Chrysene	13.99	228	1324497	44.060	ng	99
84) Bis(2-ethylhexyl)phthalate	13.94	149	115745	4.993	ng	99
86) Indeno(1,2,3-cd)pyrene	16.81	276	568699	23.436	ng	# 91
88) Benzo(b)fluoranthene	14.99	252	1759398m	59.819	ng	
89) Benzo(k)fluoranthene	15.01	252	506861m	18.329	ng	
90) Benzo(a)pyrene	15.34	252	1099099	40.580	ng	98
91) Dibenzo(a,h)anthracene	16.82	278	152734	7.179	ng	96
92) Benzo(g,h,i)perylene	17.24	276	512690	24.075	ng	95

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

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