

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103716.D
 Acq On : 17 Mar 2018 00:13
 Operator : JU/SJ
 Sample : J1940-01 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1-5

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:02:45 PM

Quant Time: Mar 19 11:14:36 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF031518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 18:31:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	156011	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	598977	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	211750	20.00	ng	0.01
63) Phenanthrene-d10	11.52	188	319848	20.00	ng	0.02
75) Chrysene-d12	14.16	240	308992	20.00	ng	0.01
86) Perylene-d12	15.69	264	245152	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	554470	54.06	ng	0.00
7) Phenol-d6	6.58	99	673273	53.65	ng	0.00
23) Nitrobenzene-d5	7.53	82	376162	43.79	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	98384	54.04	ng	0.01
44) 2-Fluorobiphenyl	9.33	172	523972	39.47	ng	0.00
78) Terphenyl-d14	13.10	244	392401	29.48	ng	0.01
Target Compounds						
22) Acetophenone	7.36	105	47485	3.085	ng	# 56
31) Naphthalene	8.27	128	62350	2.061	ng	# 72
37) 2-Methylnaphthalene	8.97	142	120585	6.284	ng	# 92
45) 1,1'-Biphenyl	9.43	154	43904	2.487	ng	95
48) Acenaphthylene	9.88	152	93145	4.283	ng	# 74
49) Dimethylphthalate	9.72	163	47295	2.928	ng	# 76
57) Fluorene	10.57	166	54222	3.789	ng	# 73
70) Phenanthrene	11.54	178	161284	9.218	ng	97
71) Anthracene	11.59	178	69373m	3.904	ng	
74) Fluoranthene	12.73	202	153257	8.502	ng	97
77) Pyrene	12.96	202	260461	11.478	ng	99
80) Benzo(a)anthracene	14.15	228	121226m	6.198	ng	
82) Chrysene	14.19	228	136067	7.298	ng	97
83) Bis(2-ethylhexyl)phthalate	14.12	149	64201	4.470	ng	93
85) Indeno(1,2,3-cd)pyrene	17.25	276	54111	3.323	ng	97
87) Benzo(b)fluoranthene	15.24	252	128091m	8.270	ng	
88) Benzo(k)fluoranthene	15.26	252	39265m	2.637	ng	
89) Benzo(a)pyrene	15.63	252	107970	7.686	ng	# 92
91) Benzo(a,h,i)perylene	17.73	276	70439	5.952	ng	97

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

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