

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103715.D
 Acq On : 16 Mar 2018 23:47
 Operator : JU/SJ
 Sample : J1938-09
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CS-3

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:02:42 PM

Quant Time: Mar 17 03:56:59 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	169407	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	668874	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	285420	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	406979	20.00	ng	0.00
75) Chrysene-d12	14.15	240	318911	20.00	ng	0.00
86) Perylene-d12	15.69	264	274790	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1062777	95.42	ng	0.01
7) Phenol-d6	6.59	99	1302126	95.56	ng	0.00
23) Nitrobenzene-d5	7.53	82	771269	80.41	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	227659	92.77	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1158801	64.76	ng	0.00
78) Terphenyl-d14	13.09	244	808274	58.83	ng	0.00
Target Compounds						
10) Phenol	6.60	94	104719	7.232	ng	95
49) Dimethylphthalate	9.72	163	152591	7.007	ng	99
70) Phenanthrene	11.53	178	62454	2.805	ng	97
74) Fluoranthene	12.72	202	201963	8.806	ng	99
77) Pyrene	12.95	202	202476	8.645	ng	98
80) Benzo(a)anthracene	14.14	228	133886	6.632	ng	96
82) Chrysene	14.18	228	129028	6.705	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	69870	4.158	ng	# 93
87) Benzo(b)fluoranthene	15.23	252	186201m	10.726	ng	
88) Benzo(k)fluoranthene	15.26	252	54562m	3.270	ng	
89) Benzo(a)pyrene	15.62	252	121688	7.728	ng	98
91) Benzo(a,h,i)perylene	17.73	276	66549	5.017	ng	96

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

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