

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF032018\
 Data File : BF103818.D
 Acq On : 20 Mar 2018 22:21
 Operator : JU/SJ
 Sample : J1966-07
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FRAC-1

Manual Integrations
 APPROVED

Sohil
 3/22/2018 12:57:52 PM

Quant Time: Mar 21 12:54:21 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	150707	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	508870	20.00	ng	0.00
38) Acenaphthene-d10	10.03	164	172304	20.00	ng	0.02
63) Phenanthrene-d10	11.53	188	133476	20.00	ng	0.03
75) Chrysene-d12	14.20	240	65639	20.00	ng	0.05
86) Perylene-d12	15.79	264	204809	20.00	ng	0.11
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	585005	59.04	ng	0.00
7) Phenol-d6	6.59	99	453669	37.42	ng	0.00
23) Nitrobenzene-d5	7.54	82	877233	120.22	ng	0.00
41) 2,4,6-Tribromophenol	10.83	330	231165	156.04	ng	0.03
44) 2-Fluorobiphenyl	9.34	172	1139284	105.47	ng	0.01
78) Terphenyl-d14	13.13	244	531769	188.05	ng	0.04
Target Compounds						
10) Phenol	6.60	94	31894	2.476	ng	97
37) 2-Methylnaphthalene	8.98	142	210897	12.937	ng	96
51) Acenaphthene	10.06	154	44496	4.235	ng	# 64
57) Fluorene	10.58	166	82846	7.114	ng	# 82
70) Phenanthrene	11.55	178	114760	15.717	ng	# 91

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

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