

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF031618\
 Data File : BF103729.D
 Acq On : 17 Mar 2018 8:33
 Operator : JU/SJ
 Sample : J1947-01
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-013-A

Manual Integrations
 APPROVED

Sohil
 3/19/2018 2:03:10 PM

Quant Time: Mar 19 08:15:14 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	148327	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	574485	20.00	ng	0.00
38) Acenaphthene-d10	10.01	164	232737	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	332184	20.00	ng	0.00
75) Chrysene-d12	14.15	240	294911	20.00	ng	0.00
86) Perylene-d12	15.69	264	236216	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1174230	120.41	ng	0.02
7) Phenol-d6	6.60	99	1427334	119.63	ng	0.01
23) Nitrobenzene-d5	7.53	82	854493	103.73	ng	0.00
41) 2,4,6-Tribromophenol	10.80	330	237673	118.77	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1307272	89.60	ng	0.00
78) Terphenyl-d14	13.09	244	928514	73.08	ng	0.00
Target Compounds						
10) Phenol	6.60	94	40228	3.173	ng	95
49) Dimethylphthalate	9.72	163	133260	7.505	ng	99
70) Phenanthrene	11.53	178	94996	5.228	ng	98
74) Fluoranthene	12.72	202	109807	5.866	ng	96
77) Pyrene	12.95	202	106674	4.925	ng	99
80) Benzo(a)anthracene	14.14	228	56387	3.021	ng	96
82) Chrysene	14.18	228	50845	2.857	ng	98
87) Benzo(b)fluoranthene	15.23	252	52925m	3.546	ng	
89) Benzo(a)pyrene	15.62	252	36077	2.665	ng	# 94

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

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