

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF033018\
 Data File : BF104115.D
 Acq On : 31 Mar 2018 1:18
 Operator : JU/SJ
 Sample : J2142-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GB-RT-4632

Manual Integrations
 APPROVED

Sohil
 4/2/2018 5:06:05 PM

Quant Time: Mar 31 04:20:42 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 30 17:20:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	111142	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	464497	20.00	ng	0.00
38) Acenaphthene-d10	9.99	164	203335	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	325738	20.00	ng	0.00
75) Chrysene-d12	14.14	240	241410	20.00	ng	0.00
86) Perylene-d12	15.68	264	210135	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	770315	110.53	ng	0.01
7) Phenol-d6	6.59	99	1019169	118.86	ng	0.00
23) Nitrobenzene-d5	7.52	82	597324	78.64	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	225798	99.73	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	945536	73.33	ng	0.00
78) Terphenyl-d14	13.07	244	732780	70.09	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	69281	4.316	ng	99
70) Phenanthrene	11.52	178	86120	4.883	ng	98
74) Fluoranthene	12.71	202	147420	7.864	ng	96
77) Pyrene	12.94	202	123868	7.138	ng	98
80) Benzo(a)anthracene	14.13	228	62380	4.130	ng	98
82) Chrysene	14.17	228	52034	3.517	ng	98
87) Benzo(b)fluoranthene	15.22	252	54618m	4.271	ng	
88) Benzo(k)fluoranthene	15.24	252	28776m	2.389	ng	
89) Benzo(a)pyrene	15.62	252	43482	3.669	ng	# 90

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

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