

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF033018\
 Data File : BF104109.D
 Acq On : 30 Mar 2018 22:39
 Operator : JU/SJ
 Sample : J1760-09 2X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-032918

Manual Integrations
 APPROVED

Sohil
 4/2/2018 5:05:54 PM

Quant Time: Mar 31 00:22:57 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	111432	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	449568	20.00	ng	0.00
38) Acenaphthene-d10	9.99	164	195859	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	304552	20.00	ng	0.00
75) Chrysene-d12	14.14	240	236382	20.00	ng	0.00
86) Perylene-d12	15.68	264	199281	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	360933	51.65	ng	0.00
7) Phenol-d6	6.59	99	484650	56.38	ng	0.00
23) Nitrobenzene-d5	7.52	82	297862	40.52	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	110436	50.64	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	561545	45.21	ng	0.00
78) Terphenyl-d14	13.07	244	409479	40.00	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.70	163	33844	2.189	ng	99
70) Phenanthrene	11.52	178	75483	4.578	ng	96
74) Fluoranthene	12.71	202	134567	7.678	ng	97
77) Pyrene	12.94	202	110759	6.518	ng	99
80) Benzo(a)anthracene	14.13	228	65510	4.429	ng	96
82) Chrysene	14.17	228	52044	3.592	ng	99
87) Benzo(b)fluoranthene	15.22	252	58989m	4.864	ng	
88) Benzo(k)fluoranthene	15.24	252	31536m	2.760	ng	
89) Benzo(a)pyrene	15.62	252	44081	3.922	ng	# 85
91) Benzo(q,h,i)perylene	17.74	276	21611	2.076	ng	# 92

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

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