

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031218\
 Data File : BF103582.D
 Acq On : 12 Mar 2018 15:20
 Operator : SJ/JU
 Sample : J1876-04
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-11-01-03

Manual Integrations
 APPROVED

Sohil
 3/13/2018 4:13:05 PM

Quant Time: Mar 12 16:24:20 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 11:47:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	136422	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	585098	20.00	ng	-0.01
38) Acenaphthene-d10	9.90	164	236851	20.00	ng	-0.01
63) Phenanthrene-d10	11.39	188	365832	20.00	ng	-0.01
75) Chrysene-d12	14.05	240	218591	20.00	ng	-0.01
86) Perylene-d12	15.56	264	204067	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	380016m	42.13	ng	0.01
7) Phenol-d6	6.50	99	972382	88.10	ng	0.00
23) Nitrobenzene-d5	7.43	82	637485	62.22	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	11023	5.50	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	952912	65.58	ng	-0.01
78) Terphenyl-d14	12.98	244	608994	66.97	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.61	163	87094	4.584	ng	99
70) Phenanthrene	11.42	178	52365	2.601	ng	97
74) Fluoranthene	12.62	202	127060	6.280	ng	97
77) Pyrene	12.85	202	131411	7.711	ng	97
80) Benzo(a)anthracene	14.04	228	58413	4.119	ng	97
82) Chrysene	14.08	228	61511m	4.467	ng	
85) Indeno(1,2,3-cd)pyrene	17.12	276	34031	3.121	ng	94
87) Benzo(b)fluoranthene	15.12	252	74898m	5.582	ng	
88) Benzo(k)fluoranthene	15.14	252	28549m	2.260	ng	
89) Benzo(a)pyrene	15.50	252	56518	4.717	ng	# 94
91) Benzo(a,h,i)perylene	17.59	276	31130	3.440	ng	# 87

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

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