

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030818\
 Data File : BF103525.D
 Acq On : 8 Mar 2018 19:09
 Operator : SJ/JU
 Sample : J1820-01
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1

Manual Integrations
 APPROVED

Sohil
 3/9/2018 11:39:33 AM

Quant Time: Mar 09 00:41:40 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 08 14:20:19 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	146699	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	620897	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	231560	20.00	ng	0.00
63) Phenanthrene-d10	11.40	188	345443	20.00	ng	0.00
75) Chrysene-d12	14.06	240	239515	20.00	ng	0.00
86) Perylene-d12	15.56	264	207474	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	979496	100.98	ng	0.00
7) Phenol-d6	6.50	99	1200169	101.12	ng	0.00
23) Nitrobenzene-d5	7.43	82	821778	75.58	ng	0.00
41) 2,4,6-Tribromophenol	10.70	330	168292	85.86	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1071548	78.20	ng	0.00
78) Terphenyl-d14	12.99	244	671091	67.35	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	11.43	178	62010	3.262	ng	98
73) Di-n-butylphthalate	11.95	149	363000	14.943	ng	99
74) Fluoranthene	12.62	202	104198	5.454	ng	99
77) Pyrene	12.85	202	97192	5.205	ng	97
80) Benzo(a)anthracene	14.04	228	47910	3.083	ng	97
82) Chrysene	14.09	228	46968	3.113	ng	99
87) Benzo(b)fluoranthene	15.12	252	60089m	4.405	ng	
89) Benzo(a)pyrene	15.50	252	39668	3.256	ng	# 93
91) Benzo(g,h,i)perylene	17.59	276	23396	2.543	ng	97

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

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