

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103339.D
 Acq On : 1 Mar 2018 16:23
 Operator : SJ/JU
 Sample : J1720-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-36

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:42 PM

Quant Time: Mar 02 09:53:02 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 17:05:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	133051	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	503316	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	185625	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	277659	20.00	ng	0.00
75) Chrysene-d12	14.06	240	183020	20.00	ng	0.00
86) Perylene-d12	15.57	264	130156	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	937370	106.55	ng	0.01
7) Phenol-d6	6.52	99	1092665	101.50	ng	0.00
23) Nitrobenzene-d5	7.44	82	680403	77.20	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	141959	90.35	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	881778	80.88	ng	0.00
78) Terphenyl-d14	12.99	244	621315	81.61	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.63	163	98173	6.593	ng	Qvalue # 99
70) Phenanthrene	11.43	178	210238	13.759	ng	98
71) Anthracene	11.49	178	71046	4.534	ng	100
74) Fluoranthene	12.63	202	401442	26.144	ng	99
77) Pyrene	12.86	202	385648	27.026	ng	99
80) Benzo(a)anthracene	14.05	228	183335	15.439	ng	97
82) Chrysene	14.09	228	166443	14.435	ng	97
85) Indeno(1,2,3-cd)pyrene	17.10	276	61601	6.748	ng	93
87) Benzo(b)fluoranthene	15.13	252	153886m	17.982	ng	
88) Benzo(k)fluoranthene	15.16	252	55702m	6.912	ng	
89) Benzo(a)pyrene	15.51	252	104642	13.693	ng	# 88
91) Benzo(a,h,i)perylene	17.57	276	59760	10.355	ng	# 89

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

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