

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF030118\
 Data File : BF103334.D
 Acq On : 1 Mar 2018 14:09
 Operator : SJ/JU
 Sample : J1720-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-44

Manual Integrations
 APPROVED

Sohil
 3/2/2018 3:30:32 PM

Quant Time: Mar 02 09:38:38 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	150839	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	617148	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	245319	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	349629	20.00	ng	0.00
75) Chrysene-d12	14.06	240	234531	20.00	ng	0.00
86) Perylene-d12	15.57	264	180584	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1055735	105.86	ng	0.01
7) Phenol-d6	6.52	99	1265730	103.72	ng	0.00
23) Nitrobenzene-d5	7.44	82	848108	78.48	ng	0.00
41) 2,4,6-Tribromophenol	10.71	330	170342	82.03	ng	0.00
44) 2-Fluorobiphenyl	9.23	172	1109188	75.92	ng	0.00
78) Terphenyl-d14	12.99	244	716363	73.42	ng	0.00
Target Compounds						
10) Phenol	6.53	94	89204	6.296	ng	94
49) Dimethylphthalate	9.62	163	127497	6.479	ng	99
70) Phenanthrene	11.43	178	290071	15.076	ng	98
71) Anthracene	11.49	178	89778	4.550	ng	99
74) Fluoranthene	12.63	202	484268	25.046	ng	99
77) Pyrene	12.86	202	474219	25.934	ng	99
80) Benzo(a)anthracene	14.05	228	228232	14.998	ng	97
82) Chrysene	14.09	228	202185	13.684	ng	98
85) Indeno(1,2,3-cd)pyrene	17.10	276	75044	6.415	ng	95
87) Benzo(b)fluoranthene	15.13	252	207844m	17.505	ng	
88) Benzo(k)fluoranthene	15.16	252	77250m	6.909	ng	
89) Benzo(a)pyrene	15.51	252	143990	13.580	ng	# 92
90) Dibenzo(a,h)anthracene	17.10	278	20498	2.502	ng	# 56
91) Benzo(a,h,i)perylene	17.57	276	74287	9.277	ng	# 87

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

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