

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031119\
 Data File : BF112959.D
 Acq On : 11 Mar 2019 17:20
 Operator : JU/SJ
 Sample : K1785-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 2019-03-05-DUP-SD

Manual Integrations
 APPROVED

Sohil
 3/12/2019 4:28:37 PM

Quant Time: Mar 12 05:02:20 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	158238	20.00	ng	-0.01
21) Naphthalene-d8	8.01	136	626793	20.00	ng	-0.01
39) Acenaphthene-d10	9.76	164	329285	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	742468	20.00	ng	0.00
76) Chrysene-d12	13.86	240	600705	20.00	ng	-0.01
87) Perylene-d12	15.27	264	555817	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1094633	107.61	ng	0.00
7) Phenol-d6	6.37	99	1422963	111.36	ng	0.00
23) Nitrobenzene-d5	7.29	82	885745	84.56	ng	-0.02
42) 2,4,6-Tribromophenol	10.55	330	476950	136.44	ng	-0.01
45) 2-Fluorobiphenyl	9.09	172	1308732	61.15	ng	-0.01
79) Terphenyl-d14	12.82	244	1570994	50.70	ng	0.00
Target Compounds						
10) Phenol	6.38	94	35938	2.410	ng	91
50) Dimethylphthalate	9.48	163	125664	5.176	ng	97
71) Phenanthrene	11.26	178	424703	11.497	ng	98
72) Anthracene	11.31	178	132918	3.437	ng	98
75) Fluoranthene	12.44	202	1334995	33.731	ng	98
78) Pyrene	12.67	202	1216024	24.013	ng	100
81) Benzo(a)anthracene	13.85	228	687429	16.512	ng	97
83) Chrysene	13.89	228	587168	14.398	ng	97
86) Indeno(1,2,3-cd)pyrene	16.62	276	288336m	8.293	ng	
88) Benzo(b)fluoranthene	14.87	252	653960m	18.190	ng	
89) Benzo(k)fluoranthene	14.89	252	262192m	8.051	ng	
90) Benzo(a)pyrene	15.21	252	486377	15.295	ng	97
91) Dibenzo(a,h)anthracene	16.63	278	81925m	3.019	ng	
92) Benzo(g,h,i)perylene	17.03	276	256664	9.420	ng	98

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

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