

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032919\
 Data File : BF113455.D
 Acq On : 30 Mar 2019 12:04
 Operator : JU/SJ
 Sample : K2153-19
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CONC-10

Manual Integrations
 APPROVED

Sohil
 4/1/2019 5:46:49 PM

Quant Time: Apr 01 05:05:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF032519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 26 03:32:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	115278	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	420798	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	193094	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	372032	20.00	ng	0.00
76) Chrysene-d12	13.84	240	381348	20.00	ng	0.00
87) Perylene-d12	15.25	264	278472	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	740361	105.01	ng	0.00
7) Phenol-d6	6.36	99	1055134	118.30	ng	0.00
23) Nitrobenzene-d5	7.27	82	621828	87.71	ng	0.00
42) 2,4,6-Tribromophenol	10.53	330	200060	83.88	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	1197920	100.44	ng	0.00
79) Terphenyl-d14	12.80	244	1337021	77.28	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.46	163	58359	4.078	ng	98
71) Phenanthrene	11.24	178	109826	5.945	ng	98
75) Fluoranthene	12.42	202	179404	8.590	ng	99
78) Pyrene	12.65	202	194832	7.373	ng	98
81) Benzo(a)anthracene	13.83	228	84267	3.860	ng	98
83) Chrysene	13.87	228	75233	3.393	ng	99
88) Benzo(b)fluoranthene	14.85	252	68921m	3.984	ng	
90) Benzo(a)pyrene	15.19	252	51962	3.388	ng	# 95
92) Benzo(g,h,i)perylene	17.02	276	28458	2.128	ng	# 90

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

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