

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031020\
 Data File : BF119319.D
 Acq On : 10 Mar 2020 16:22
 Operator : CG/JU
 Sample : L1852-19
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3-D

Manual Integrations
 APPROVED

mohammad
 3/11/2020 4:58:01 PM

Quant Time: Mar 11 03:51:46 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 06 23:38:46 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	104370	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	390378	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	191378	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	298620	20.00	ng	0.00
76) Chrysene-d12	13.89	240	238629	20.00	ng	0.00
87) Perylene-d12	15.32	264	223606	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	745537	119.38	ng	0.01
7) Phenol-d6	6.39	99	905146	119.17	ng	0.00
23) Nitrobenzene-d5	7.29	82	589263	79.70	ng	0.00
42) 2,4,6-Tribromophenol	10.55	330	236919	94.63	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	1110011	85.45	ng	0.00
79) Terphenyl-d14	12.83	244	1029445	74.27	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.48	163	160600	10.974	ng	Qvalue 97
75) Fluoranthene	12.46	202	74079	4.285	ng	99
78) Pyrene	12.69	202	72415	3.779	ng	98
81) Benzo(a)anthracene	13.87	228	43057	2.648	ng	97
83) Chrysene	13.92	228	42392	2.617	ng	96
88) Benzo(b)fluoranthene	14.92	252	51659m	3.505	ng	
90) Benzo(a)pyrene	15.26	252	30787	2.314	ng	# 95

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

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