

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF033020\
 Data File : BF119639.D
 Acq On : 30 Mar 2020 18:05
 Operator : MA/SJ
 Sample : L1762-23 5X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-032720

Manual Integrations
 APPROVED

mohammad
 3/31/2020 3:16:37 PM

Quant Time: Mar 31 00:51:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 27 07:52:09 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	306329	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1102129	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	508948	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	829324	20.00	ng	0.00
76) Chrysene-d12	14.02	240	563578	20.00	ng	0.00
87) Perylene-d12	15.48	264	531842	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	359353	18.67	ng	0.00
7) Phenol-d6	6.45	99	494792	20.13	ng	-0.02
23) Nitrobenzene-d5	7.39	82	222503	10.97	ng	-0.02
42) 2,4,6-Tribromophenol	10.66	330	66708	12.70	ng	-0.01
45) 2-Fluorobiphenyl	9.20	172	587986	17.12	ng	0.00
79) Terphenyl-d14	12.96	244	485179	16.87	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.59	163	79410	2.197	ng	99
71) Phenanthrene	11.39	178	882141	20.514	ng	98
72) Anthracene	11.44	178	191157	4.374	ng	97
75) Fluoranthene	12.59	202	1333369	28.651	ng	98
78) Pyrene	12.82	202	1448617	31.320	ng	98
81) Benzo(a)anthracene	14.00	228	604943	16.507	ng	96
83) Chrysene	14.04	228	660786	17.471	ng	98
86) Indeno(1,2,3-cd)pyrene	16.93	276	252019	7.075	ng	98
88) Benzo(b)fluoranthene	15.05	252	639954m	18.013	ng	
89) Benzo(k)fluoranthene	15.07	252	213705m	6.317	ng	
90) Benzo(a)pyrene	15.42	252	493829	15.295	ng	98
91) Dibenzo(a,h)anthracene	16.96	278	69203	2.451	ng	94
92) Benzo(a,h,i)perylene	17.38	276	245521	8.654	ng	93

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

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