

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042420\
 Data File : BF120201.D
 Acq On : 24 Apr 2020 14:45
 Operator : MA/SJ
 Sample : L2377-01 10X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MILL-1

Manual Integrations
 APPROVED

mohammad
 4/27/2020 3:04:38 PM

Quant Time: Apr 24 16:09:41 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 24 10:24:30 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	130340	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	495684	20.00	ng	0.00
39) Acenaphthene-d10	9.70	164	232444	20.00	ng	0.00
64) Phenanthrene-d10	11.18	188	378913	20.00	ng	0.00
76) Chrysene-d12	13.82	240	248504	20.00	ng	-0.01
87) Perylene-d12	15.22	264	231346	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.27	112	49699	6.59	ng	0.00
7) Phenol-d6	6.32	99	65725	6.76	ng	0.00
23) Nitrobenzene-d5	7.22	82	42680	4.45	ng	-0.02
42) 2,4,6-Tribromophenol	10.49	330	13674	4.80	ng	-0.01
45) 2-Fluorobiphenyl	9.02	172	89751	5.48	ng	-0.01
79) Terphenyl-d14	12.76	244	73244	4.96	ng	-0.01
Target Compounds						
71) Phenanthrene	11.20	178	80412	3.987	ng	96
75) Fluoranthene	12.39	202	78537	3.609	ng	96
78) Pyrene	12.62	202	87320	4.168	ng	98
81) Benzo(a)anthracene	13.80	228	41745m	2.553	ng	
83) Chrysene	13.84	228	39696	2.390	ng	94
88) Benzo(b)fluoranthene	14.82	252	40179m	2.414	ng	
90) Benzo(a)pyrene	15.16	252	33869	2.420	ng	98

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

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