

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF060420\
 Data File : BF120551.D
 Acq On : 4 Jun 2020 23:29
 Operator : JU/CG
 Sample : L2858-01 10X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHRT27247

Manual Integrations
 APPROVED

mohammad
 6/5/2020 11:58:22 AM

Quant Time: Jun 05 01:55:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF052820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 04 16:38:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	166265	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	663656	20.00	ng	0.00
39) Acenaphthene-d10	9.87	164	248729	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	443323	20.00	ng	0.00
76) Chrysene-d12	14.00	240	306679	20.00	ng	0.00
86) Perylene-d12	15.46	264	371629	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	73220	8.39	ng	0.00
7) Phenol-d6	6.46	99	89053	8.28	ng	-0.01
23) Nitrobenzene-d5	7.38	82	51786	5.10	ng	-0.02
42) 2,4,6-Tribromophenol	10.65	330	20423	8.86	ng	-0.01
45) 2-Fluorobiphenyl	9.18	172	96507	6.58	ng	-0.01
79) Terphenyl-d14	12.94	244	101352	7.35	ng	0.00
Target Compounds						
75) Fluoranthene	12.57	202	55652	2.648	ng	98
78) Pyrene	12.80	202	109802	5.148	ng	99
81) Benzo(a)anthracene	13.99	228	37026m	2.253	ng	
83) Chrysene	14.02	228	42682m	2.477	ng	
88) Benzo(b)fluoranthene	15.03	252	58159m	2.818	ng	
90) Benzo(a)pyrene	15.39	252	42020	2.327	ng	# 87

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

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