

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081720\
 Data File : BF121314.D
 Acq On : 17 Aug 2020 17:20
 Operator : JU/CG
 Sample : L3508-19 2X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-08-081420

Manual Integrations
 APPROVED

mohammad
 8/18/2020 1:50:18 PM

Quant Time: Aug 17 17:54:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 10:52:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	91630	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	362953	20.00	ng	0.00
39) Acenaphthene-d10	9.74	164	183549	20.00	ng	0.00
64) Phenanthrene-d10	11.22	188	344779	20.00	ng	0.00
76) Chrysene-d12	13.86	240	261258	20.00	ng	0.00
86) Perylene-d12	15.27	264	286280	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	62167	10.29	ng	0.00
7) Phenol-d6	6.34	99	95723	12.29	ng	0.00
23) Nitrobenzene-d5	7.26	82	69272	10.43	ng	-0.01
42) 2,4,6-Tribromophenol	10.53	330	16167	7.15	ng	0.00
45) 2-Fluorobiphenyl	9.06	172	150571	11.35	ng	0.00
79) Terphenyl-d14	12.80	244	146789	10.95	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.46	163	32289	2.338	ng	98
78) Pyrene	12.66	202	45856	2.357	ng	98
81) Benzo(a)anthracene	13.84	228	43242	2.440	ng	# 87
83) Chrysene	13.88	228	36986	2.044	ng	97
88) Benzo(b)fluoranthene	14.87	252	64988m	3.388	ng	
90) Benzo(a)pyrene	15.21	252	43553	2.516	ng	# 92

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

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