

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042020\
 Data File : BF120091.D
 Acq On : 20 Apr 2020 21:24
 Operator : MA/SJ
 Sample : L2314-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WB-4-1

Manual Integrations
 APPROVED

mohammad
 4/21/2020 2:55:09 PM

Quant Time: Apr 21 03:57:17 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 20 14:10:57 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	154579	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	572596	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	278348	20.00	ng	0.00
64) Phenanthrene-d10	11.24	188	478816	20.00	ng	-0.01
76) Chrysene-d12	13.89	240	320397	20.00	ng	-0.01
87) Perylene-d12	15.31	264	286058	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	637229	71.24	ng	0.00
7) Phenol-d6	6.36	99	824973	71.58	ng	0.00
23) Nitrobenzene-d5	7.28	82	512784	46.33	ng	-0.01
42) 2,4,6-Tribromophenol	10.55	330	195613	57.40	ng	-0.01
45) 2-Fluorobiphenyl	9.08	172	984605	50.22	ng	0.00
79) Terphenyl-d14	12.83	244	915856	48.10	ng	-0.01
Target Compounds						
31) Naphthalene	8.02	128	135806	4.508	ng	99
50) Dimethylphthalate	9.47	163	181925	8.450	ng	99
71) Phenanthrene	11.26	178	93634	3.674	ng	97
75) Fluoranthene	12.46	202	87802	3.193	ng	96
78) Pyrene	12.68	202	110636	4.096	ng	99
81) Benzo(a)anthracene	13.87	228	64983	3.083	ng	# 84
83) Chrysene	13.91	228	67983	3.174	ng	# 91
86) Indeno(1,2,3-cd)pyrene	16.69	276	65451	3.467	ng	97
88) Benzo(b)fluoranthene	14.90	252	101797m	4.947	ng	
90) Benzo(a)pyrene	15.25	252	83030	4.799	ng	# 93
92) Benzo(g,h,i)perylene	17.11	276	71549	4.729	ng	97

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

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