

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061820\
 Data File : BF120651.D
 Acq On : 18 Jun 2020 19:33
 Operator : JU/CG
 Sample : L3037-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SL-WC-01C

Manual Integrations
 APPROVED

mohammad
 6/19/2020 12:58:04 PM

Quant Time: Jun 19 01:33:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 11 11:28:22 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	153767	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	563384	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	283536	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	530935	20.00	ng	0.00
76) Chrysene-d12	13.96	240	436136	20.00	ng	0.00
86) Perylene-d12	15.40	264	388577	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	981839	109.47	ng	0.00
7) Phenol-d6	6.44	99	1218393	109.07	ng	0.00
23) Nitrobenzene-d5	7.36	82	749016	74.31	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	354908	107.66	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1446585	78.06	ng	0.00
79) Terphenyl-d14	12.91	244	1632473	82.68	ng	0.00
Target Compounds						
49) Acenaphthylene	9.70	152	60138	2.251	ng	98
50) Dimethylphthalate	9.55	163	137875	6.752	ng	98
71) Phenanthrene	11.35	178	237091	8.983	ng	98
72) Anthracene	11.40	178	56864	2.137	ng	99
75) Fluoranthene	12.53	202	280077	9.282	ng	98
78) Pyrene	12.76	202	383610	12.416	ng	98
81) Benzo(a)anthracene	13.95	228	182074m	6.949	ng	
83) Chrysene	13.99	228	187725	6.984	ng	97
87) Indeno(1,2,3-cd)pyrene	16.83	276	65491	2.771	ng	96
88) Benzo(b)fluoranthene	14.99	252	166179m	7.179	ng	
89) Benzo(k)fluoranthene	15.01	252	51696m	2.524	ng	
90) Benzo(a)pyrene	15.34	252	124516	6.011	ng	# 93
92) Benzo(a,h,i)perylene	17.26	276	77505	4.057	ng	93

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

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