

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062220\
 Data File : BF120692.D
 Acq On : 22 Jun 2020 23:04
 Operator : JU/CG
 Sample : L3082-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-18

Manual Integrations
 APPROVED

Sohil
 6/23/2020 2:03:45 PM

Quant Time: Jun 23 03:12:28 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	145793	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	550896	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	285475	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	576722	20.00	ng	0.00
76) Chrysene-d12	13.98	240	356223	20.00	ng	0.02
86) Perylene-d12	15.43	264	448223	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	791795	93.11	ng	0.00
7) Phenol-d6	6.43	99	1016981	96.02	ng	0.00
23) Nitrobenzene-d5	7.36	82	640135	64.95	ng	-0.01
42) 2,4,6-Tribromophenol	10.63	330	321739	96.94	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1239511	66.43	ng	0.00
79) Terphenyl-d14	12.92	244	1528686	94.79	ng	0.00
Target Compounds						
31) Naphthalene	8.10	128	215657	7.972	ng	99
37) 2-Methylnaphthalene	8.79	142	61898	3.505	ng	100
38) 1-Methylnaphthalene	8.89	142	60675	3.651	ng	98
49) Acenaphthylene	9.69	152	563919	20.962	ng	98
50) Dimethylphthalate	9.55	163	113778	5.534	ng	99
52) Acenaphthene	9.86	154	120750	7.526	ng	100
55) Dibenzofuran	10.04	168	276882	11.255	ng	98
58) Fluorene	10.38	166	348733	18.388	ng	99
71) Phenanthrene	11.36	178	3583865	125.010	ng	98
72) Anthracene	11.40	178	855172	29.593	ng	100
73) Carbazole	11.55	167	527487	18.440	ng	99
75) Fluoranthene	12.56	202	5110538	155.915	ng	98
78) Pyrene	12.79	202	4742449	187.928	ng	95
81) Benzo(a)anthracene	13.97	228	2700235	126.170	ng	# 90
83) Chrysene	14.01	228	2464124m	112.247	ng	
87) Indeno(1,2,3-cd)pyrene	16.89	276	2061070	75.595	ng	96
88) Benzo(b)fluoranthene	15.02	252	3289572m	123.207	ng	
89) Benzo(k)fluoranthene	15.04	252	1111135m	47.031	ng	
90) Benzo(a)pyrene	15.38	252	2502151	104.714	ng	96
91) Dibenz(a,h)anthracene	16.89	278	465842m	21.307	ng	
92) Benzo(g,h,i)perylene	17.32	276	1998198	90.672	ng	100

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

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