

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011620\
 Data File : BP001649.D
 Acq On : 17 Jan 2020 03:07
 Operator : JU
 Sample : L1108-06 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 INWOOD-BH-03-B

Manual Integrations
 APPROVED

mohammad
 1/17/2020 2:09:03 PM

Quant Time: Jan 17 06:32:26 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 14 21:52:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	37768	20.00	ng	0.00
21) Naphthalene-d8	10.41	136	145179	20.00	ng	0.00
39) Acenaphthene-d10	14.28	164	98176	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	213482	20.00	ng	0.00
76) Chrysene-d12	21.15	240	162718	20.00	ng	0.00
87) Perylene-d12	23.39	264	173350	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	44294	22.97	ng	0.00
7) Phenol-d6	6.83	99	66592	21.61	ng	0.00
23) Nitrobenzene-d5	8.78	82	43202	13.00	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	22696	15.19	ng	0.00
45) 2-Fluorobiphenyl	12.90	172	97099	14.54	ng	0.00
79) Terphenyl-d14	19.63	244	136614	15.36	ng	-0.01
Target Compounds						
49) Acenaphthylene	14.00	152	106243	11.782	ng	99
71) Phenanthrene	17.09	178	97729	8.418	ng	99
72) Anthracene	17.18	178	113111	10.008	ng	98
75) Fluoranthene	19.07	202	869957	56.947	ng	100
78) Pyrene	19.43	202	799236	66.865	ng	99
81) Benzo(a)anthracene	21.14	228	477073	42.099	ng	94
83) Chrysene	21.19	228	428516	38.801	ng	96
86) Indeno(1,2,3-cd)pyrene	25.65	276	333187	25.773	ng	99
88) Benzo(b)fluoranthene	22.73	252	705118	64.550	ng	99
89) Benzo(k)fluoranthene	22.76	252	205788m	19.322	ng	
90) Benzo(a)pyrene	23.29	252	499894	48.198	ng	97
91) Dibenzo(a,h)anthracene	25.65	278	77889	7.279	ng	# 90
92) Benzo(a,h,i)perylene	26.34	276	303080	28.490	ng	97

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

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