

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP090220\
 Data File : BP003193.D
 Acq On : 02 Sep 2020 18:45
 Operator : CG/JU
 Sample : L3866-01 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-03-090120

Manual Integrations
 APPROVED

mohammad
 9/3/2020 1:48:42 PM

Quant Time: Sep 03 01:31:44 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP082420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 01 14:37:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	178578	20.00	ng	0.00
21) Naphthalene-d8	10.43	136	772742	20.00	ng	0.00
39) Acenaphthene-d10	14.30	164	512876	20.00	ng	0.00
64) Phenanthrene-d10	17.05	188	1218112	20.00	ng	0.00
76) Chrysene-d12	21.15	240	1350843	20.00	ng	0.00
86) Perylene-d12	23.38	264	1402856	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.26	112	387925	38.41	ng	0.00
7) Phenol-d6	6.84	99	486411	38.42	ng	0.00
23) Nitrobenzene-d5	8.80	82	286412	26.72	ng	0.00
42) 2,4,6-Tribromophenol	15.79	330	289707	41.77	ng	0.00
45) 2-Fluorobiphenyl	12.92	172	908735	25.95	ng	0.00
79) Terphenyl-d14	19.63	244	1678690	27.34	ng	0.00
Target Compounds						
50) Dimethylphthalate	13.76	163	159859	4.087	ng	99
71) Phenanthrene	17.09	178	579668	8.782	ng	100
72) Anthracene	17.18	178	239810	3.796	ng	99
75) Fluoranthene	19.07	202	1762071	23.098	ng	99
78) Pyrene	19.42	202	1521055	17.349	ng	100
81) Benzo(a)anthracene	21.13	228	929812	10.989	ng	98
83) Chrysene	21.18	228	785818	9.700	ng	98
87) Indeno(1,2,3-cd)pyrene	25.65	276	624802	5.972	ng	96
88) Benzo(b)fluoranthene	22.71	252	1188830	13.635	ng	96
89) Benzo(k)fluoranthene	22.75	252	382393m	4.590	ng	
90) Benzo(a)pyrene	23.28	252	887711	10.974	ng	97
92) Benzo(a,h,i)perylene	26.34	276	587244	6.809	ng	98

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

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