

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101719\
 Data File : BN008217.D
 Acq On : 17 Oct 2019 11:59
 Operator : JU
 Sample : K5175-18DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP90DL

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:08:10 PM

Quant Time: Oct 18 02:19:14 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 17 01:48:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2741	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	13096	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8548	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	18063m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	15303	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	11411	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	417	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	1287	0.02	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.96	152	1407	0.035	ng/ul#	88
9) Fluorene	15.30	166	798	0.022	ng/ul#	94
12) Phenanthrene	17.03	178	16711	0.316	ng/ul	100
13) Anthracene	17.12	178	2960	0.064	ng/ul#	94
15) Fluoranthene	19.06	202	43319	0.620	ng/ul	99
17) Pyrene	19.42	202	45127	0.741	ng/ul	100
18) Benzo(a)anthracene	21.18	228	21167	0.380	ng/ul	96
19) Chrysene	21.23	228	25293	0.370	ng/ul	98
21) Benzo(b)fluoranthene	22.73	252	23572m	0.635	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	6642m	0.158	ng/ul	
23) Benzo(a)pyrene	23.28	252	15202	0.389	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.54	276	9488	0.206	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.54	278	2129	0.059	ng/ul#	58
26) Benzo(g,h,i)perylene	26.20	276	8177	0.194	ng/ul	94

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

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