

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101719\
 Data File : BN008239.D
 Acq On : 18 Oct 2019 01:57
 Operator : JU
 Sample : K5177-19
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPC8

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:09:50 PM

Quant Time: Oct 18 04:10:05 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	2969	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	14413	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8717	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	18900m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	14705	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	11908	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	3355	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	9740	0.15	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.95	152	1133	0.027	ng/ul#	82
8) Acenaphthene	14.30	153	2329	0.072	ng/ul	99
9) Fluorene	15.29	166	3147	0.086	ng/ul	98
12) Phenanthrene	17.03	178	116395	2.103	ng/ul	99
13) Anthracene	17.12	178	13353	0.274	ng/ul	99
15) Fluoranthene	19.05	202	329384	4.507	ng/ul	99
17) Pyrene	19.41	202	246808	4.215	ng/ul	99
18) Benzo(a)anthracene	21.17	228	96767	1.810	ng/ul	99
19) Chrysene	21.22	228	114070	1.739	ng/ul	99
21) Benzo(b)fluoranthene	22.72	252	123204m	3.178	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	39824m	0.907	ng/ul	
23) Benzo(a)pyrene	23.27	252	74806	1.834	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.52	276	57389	1.195	ng/ul	96
25) Dibenzo(a,h)anthracene	25.53	278	10989	0.289	ng/ul	94
26) Benzo(g,h,i)perylene	26.18	276	51533	1.170	ng/ul	99

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

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