

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101719\
 Data File : BN008223.D
 Acq On : 17 Oct 2019 15:35
 Operator : JU
 Sample : K5177-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPG7

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:08:54 PM

Quant Time: Oct 18 02:51:00 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	7.61	152	361	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	1438	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	781	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	1394m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	1223	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	1475	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	533	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	1120	0.23	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.96	152	132	0.035	ng/ul#	10
12) Phenanthrene	17.03	178	391	0.096	ng/ul#	58
13) Anthracene	17.12	178	89	0.025	ng/ul#	1
15) Fluoranthene	19.06	202	823	0.153	ng/ul	72
17) Pyrene	19.42	202	732m	0.150	ng/ul	
18) Benzo(a)anthracene	21.18	228	477	0.107	ng/ul#	53
19) Chrysene	21.23	228	665	0.122	ng/ul#	66
21) Benzo(b)fluoranthene	22.73	252	674m	0.140	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	260m	0.048	ng/ul	
23) Benzo(a)pyrene	23.28	252	434	0.086	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.54	276	252	0.042	ng/ul#	57
26) Benzo(g,h,i)perylene	26.20	276	270	0.049	ng/ul#	1

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

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