

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101719\
 Data File : BN008227.D
 Acq On : 17 Oct 2019 18:46
 Operator : JU
 Sample : K5177-09
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP98

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:08:35 PM

Quant Time: Oct 18 02:58:39 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	2571	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	12462	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	8014	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	17909m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	16222	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	12450	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	2869	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	8950	0.14	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.42	128	784	0.022	ng/ul#	74
7) Acenaphthylene	13.95	152	1861	0.049	ng/ul#	65
12) Phenanthrene	17.03	178	12328	0.235	ng/ul	99
13) Anthracene	17.12	178	2534	0.055	ng/ul#	92
15) Fluoranthene	19.05	202	37086	0.536	ng/ul	99
17) Pyrene	19.41	202	35915	0.556	ng/ul	96
18) Benzo(a)anthracene	21.17	228	17759	0.301	ng/ul	96
19) Chrysene	21.23	228	22350	0.309	ng/ul	98
21) Benzo(b)fluoranthene	22.72	252	24142m	0.596	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	8058m	0.176	ng/ul	
23) Benzo(a)pyrene	23.28	252	15041	0.353	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.54	276	10224	0.204	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.54	278	2253	0.057	ng/ul#	62
26) Benzo(g,h,i)perylene	26.19	276	9715	0.211	ng/ul	96

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

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