

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008174.D
 Acq On : 15 Oct 2019 15:39
 Operator : JU
 Sample : K5175-06
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP86

Manual Integrations
 APPROVED

mohammad
 10/16/2019 4:08:00 PM

Quant Time: Oct 15 17:51:33 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 : Tue Oct 15 15:13:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1936	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	9795	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	6277	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	14455m	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	14022m	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	14756	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3316	0.20	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	11478m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.04	178	3738	0.088	ng/ul	96
15) Fluoranthene	19.06	202	9838m	0.176	ng/ul	
17) Pyrene	19.42	202	8216m	0.147	ng/ul	
18) Benzo(a)anthracene	21.18	228	5009	0.098	ng/ul	93
19) Chrysene	21.24	228	5054	0.081	ng/ul	95
21) Benzo(b)fluoranthene	22.73	252	6353m	0.132	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	2567m	0.047	ng/ul	
23) Benzo(a)pyrene	23.28	252	3788	0.075	ng/ul#	52
24) Indeno(1,2,3-cd)pyrene	25.55	276	2352	0.040	ng/ul#	92
26) Benzo(g,h,i)perylene	26.21	276	1878	0.034	ng/ul#	67

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

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