

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101419\
 Data File : BN008177.D
 Acq On : 15 Oct 2019 17:26
 Operator : JU
 Sample : K5175-07
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP87

Manual Integrations
 APPROVED

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

Quant Time: Oct 16 03:56:44 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	1956	0.40	ng/ul	0.00
2) Naphthalene-d8	10.38	136	9764	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	6169	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	14096m	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	14789m	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	15349	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3827	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	12592m	0.26	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.04	178	3441	0.083	ng/ul	96
15) Fluoranthene	19.06	202	10771m	0.198	ng/ul	
17) Pyrene	19.42	202	9091m	0.154	ng/ul	
18) Benzo(a)anthracene	21.18	228	6010	0.112	ng/ul	95
19) Chrysene	21.23	228	6244	0.095	ng/ul	96
21) Benzo(b)fluoranthene	22.73	252	10115m	0.202	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	3074m	0.054	ng/ul	
23) Benzo(a)pyrene	23.29	252	6210	0.118	ng/ul#	73
24) Indeno(1,2,3-cd)pyrene	25.55	276	4976	0.080	ng/ul	98
25) Dibenzo(a,h)anthracene	25.56	278	1155	0.024	ng/ul#	30
26) Benzo(g,h,i)perylene	26.21	276	4313	0.076	ng/ul#	88

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

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