

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN007018.D
 Acq On : 08 Aug 2019 10:27
 Operator : HP/JU
 Sample : K3962-06
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP28

Manual Integrations
 APPROVED

mohammad
 8/9/2019 6:16:32 AM

Quant Time: Aug 08 12:00:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	8029	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	31908	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.11	164	19483	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	49992m	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	45509m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	37641m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	8315	0.15	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	30912m	0.19	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	16.90	178	8346	0.052	ng/ul	98
15) Fluoranthene	18.93	202	24655m	0.124	ng/ul	
17) Pyrene	19.30	202	25671	0.126	ng/ul	99
18) Benzo(a)anthracene	21.06	228	11460	0.057	ng/ul	93
19) Chrysene	21.11	228	13961m	0.075	ng/ul	
21) Benzo(b)fluoranthene	22.58	252	15398m	0.101	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	5754m	0.039	ng/ul	
23) Benzo(a)pyrene	23.12	252	9972m	0.069	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.30	276	6286m	0.041	ng/ul	
26) Benzo(g,h,i)perylene	25.94	276	6390m	0.049	ng/ul	

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

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