

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008282.D
 Acq On : 21 Oct 2019 14:36
 Operator : JU
 Sample : K5199-14
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP81

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:27:15 PM

Quant Time: Oct 21 16:05:19 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 : Mon Oct 21 15:18:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1895	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	9052	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5613	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	12463m	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	12008	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	14100	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	4024	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	11637	0.27	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.94	152	544	0.020	ng/ul#	66
12) Phenanthrene	17.01	178	5205	0.143	ng/ul	98
13) Anthracene	17.10	178	1037	0.032	ng/ul#	82
15) Fluoranthene	19.04	202	11834	0.246	ng/ul	95
17) Pyrene	19.40	202	11040	0.231	ng/ul#	91
18) Benzo(a)anthracene	21.16	228	4689	0.107	ng/ul	95
19) Chrysene	21.22	228	5973	0.111	ng/ul	96
21) Benzo(b)fluoranthene	22.71	252	7391m	0.161	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	2843m	0.055	ng/ul	
23) Benzo(a)pyrene	23.26	252	4651	0.096	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	25.51	276	2788	0.049	ng/ul#	83
26) Benzo(g,h,i)perylene	26.16	276	2513	0.048	ng/ul#	80

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

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