

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008286.D
 Acq On : 21 Oct 2019 17:03
 Operator : JU
 Sample : K5178-04DL 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPE7DL

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:27:26 PM

Quant Time: Oct 22 00:09:17 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 16:03:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1804	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	8229	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5260	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	11240	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	11196	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	9558	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	807	0.06	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	2887	0.07	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.94	152	998	0.040	ng/ul#	85
12) Phenanthrene	17.01	178	7563	0.230	ng/ul	98
13) Anthracene	17.11	178	1813	0.063	ng/ul#	90
15) Fluoranthene	19.04	202	21089	0.485	ng/ul	99
17) Pyrene	19.40	202	22262	0.499	ng/ul	98
18) Benzo(a)anthracene	21.17	228	12608	0.310	ng/ul	95
19) Chrysene	21.22	228	13123	0.263	ng/ul	98
21) Benzo(b)fluoranthene	22.71	252	16538m	0.532	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	5110m	0.145	ng/ul	
23) Benzo(a)pyrene	23.26	252	10756	0.329	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.51	276	6719	0.174	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.51	278	1446	0.047	ng/ul#	58
26) Benzo(g,h,i)perylene	26.17	276	7096	0.201	ng/ul	95

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

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