

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101819\
 Data File : BN008246.D
 Acq On : 18 Oct 2019 14:25
 Operator : JU
 Sample : K5177-16
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPK2

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:40:13 AM

Quant Time: Oct 18 17:33:24 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 : Fri Oct 18 17:15:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2826	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	13289	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	8286	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	17755	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	17074	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	20519	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	1682	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	4818	0.08	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.02	178	1310	0.025	ng/ul#	86
15) Fluoranthene	19.05	202	2928	0.043	ng/ul	87
17) Pyrene	19.41	202	3265m	0.048	ng/ul	
18) Benzo(a)anthracene	21.17	228	1813	0.029	ng/ul#	83
19) Chrysene	21.22	228	1942	0.025	ng/ul#	90
21) Benzo(b)fluoranthene	22.72	252	2749m	0.041	ng/ul	
23) Benzo(a)pyrene	23.27	252	1757	0.025	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.53	276	1828	0.022	ng/ul#	77
26) Benzo(g,h,i)perylene	26.19	276	1610	0.021	ng/ul#	65

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

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