

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061819\
 Data File : BN006385.D
 Acq On : 18 Jun 2019 20:08
 Operator : JU/SJ
 Sample : K3335-04
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Z4

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:39:22 AM

Quant Time: Jun 19 02:06:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4103	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	15616	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	7805	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	14588m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10738m	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	15706	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	3435	0.14	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	6930m	0.18	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	2473	0.057	ng/ul#	91
9) Fluorene	15.35	166	1899	0.054	ng/ul#	94
12) Phenanthrene	17.09	178	38073	0.813	ng/ul	99
13) Anthracene	17.17	178	3177m	0.071	ng/ul	
15) Fluoranthene	19.11	202	52341m	1.012	ng/ul	
17) Pyrene	19.48	202	44089	0.779	ng/ul	100
18) Benzo(a)anthracene	21.24	228	16632	0.350	ng/ul	99
19) Chrysene	21.30	228	22901	0.514	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	27838m	0.425	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	9228m	0.145	ng/ul	
23) Benzo(a)pyrene	23.39	252	16120	0.264	ng/ul#	94
24) Indeno(1,2,3-cd)pyrene	25.73	276	11989	0.182	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.73	278	2830	0.054	ng/ul#	76
26) Benzo(g,h,i)perylene	26.41	276	9185	0.166	ng/ul	100

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

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