

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
 Data File : BN006243.D
 Acq On : 11 Jun 2019 10:18
 Operator : JU/SJ
 Sample : K3017-21DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51F2DL

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:39:03 PM

Quant Time: Jun 11 12:16:56 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 : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4919	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	18725	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8780	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16397	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	12473	0.40	ng/ul	0.00
20) Perylene-d12	23.51	264	16127	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1009	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	1756	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.47	128	1046	0.020	ng/ul#	75
7) Acenaphthylene	14.01	152	4186	0.086	ng/ul#	95
12) Phenanthrene	17.09	178	48841	0.928	ng/ul	99
13) Anthracene	17.19	178	5159	0.103	ng/ul#	94
15) Fluoranthene	19.12	202	122298	2.104	ng/ul	98
17) Pyrene	19.49	202	87294	1.329	ng/ul	99
18) Benzo(a)anthracene	21.26	228	11722	0.212	ng/ul	95
19) Chrysene	21.30	228	47330	0.914	ng/ul	99
21) Benzo(b)fluoranthene	22.84	252	52416m	0.779	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	17591m	0.269	ng/ul	
23) Benzo(a)pyrene	23.41	252	16528	0.264	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.74	276	16837	0.248	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.74	278	3133	0.058	ng/ul#	77
26) Benzo(g,h,i)perylene	26.42	276	15148	0.267	ng/ul	97

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

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