

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006193.D
 Acq On : 08 Jun 2019 16:13
 Operator : JU/SJ
 Sample : K3016-05
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51C4

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:30:51 PM

Quant Time: Jun 09 04:32:53 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 04:20:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4536	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	17257	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9117	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19041m	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	15796m	0.40	ng/ul	0.00
20) Perylene-d12	23.52	264	14315m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	4498	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10573m	0.21	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	1724	0.034	ng/ul#	86
12) Phenanthrene	17.09	178	4774	0.078	ng/ul	96
13) Anthracene	17.18	178	1205m	0.021	ng/ul	
15) Fluoranthene	19.12	202	10976m	0.163	ng/ul	
17) Pyrene	19.49	202	11413	0.137	ng/ul	97
18) Benzo(a)anthracene	21.26	228	4653m	0.067	ng/ul	
19) Chrysene	21.31	228	6229m	0.095	ng/ul	
21) Benzo(b)fluoranthene	22.85	252	8435m	0.141	ng/ul	
22) Benzo(k)fluoranthene	22.89	252	3338m	0.057	ng/ul	
23) Benzo(a)pyrene	23.42	252	4862	0.087	ng/ul#	60
24) Indeno(1,2,3-cd)pyrene	25.76	276	3469	0.058	ng/ul	92
25) Dibenzo(a,h)anthracene	25.76	278	973	0.020	ng/ul#	34
26) Benzo(g,h,i)perylene	26.44	276	2704	0.054	ng/ul#	81

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

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