

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060819\
 Data File : BN006192.D
 Acq On : 08 Jun 2019 15:38
 Operator : JU/SJ
 Sample : K3016-01
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51C0

Manual Integrations
 APPROVED

mohammad
 6/11/2019 4:30:46 PM

Quant Time: Jun 08 16:28:44 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 : Fri Jun 07 01:31:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	4110	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	15731	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	8153	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	16870m	0.40	ng/ul	0.00
16) Chrysene-d12	21.28	240	13474m	0.40	ng/ul	0.01
20) Perylene-d12	23.52	264	19132	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	3670	0.15	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	7639m	0.17	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	2192	0.048	ng/ul#	88
12) Phenanthrene	17.09	178	4434	0.082	ng/ul	95
13) Anthracene	17.19	178	1318m	0.025	ng/ul	
15) Fluoranthene	19.12	202	10656m	0.178	ng/ul	
17) Pyrene	19.49	202	11445	0.161	ng/ul	98
18) Benzo(a)anthracene	21.26	228	4889	0.082	ng/ul	91
19) Chrysene	21.31	228	7263	0.130	ng/ul	95
21) Benzo(b)fluoranthene	22.85	252	9418m	0.118	ng/ul	
22) Benzo(k)fluoranthene	22.89	252	3614m	0.047	ng/ul	
23) Benzo(a)pyrene	23.42	252	4358	0.059	ng/ul#	68
24) Indeno(1,2,3-cd)pyrene	25.76	276	3601	0.045	ng/ul#	91
26) Benzo(g,h,i)perylene	26.44	276	2095	0.031	ng/ul#	77

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

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