

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061219\
 Data File : BN006292.D
 Acq On : 12 Jun 2019 21:55
 Operator : JU/SJ
 Sample : K3231-13
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DB8M0

Manual Integrations
 APPROVED

mohammad
 6/14/2019 8:34:17 AM

Quant Time: Jun 13 03:08:08 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 : Wed Jun 12 17:24:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	5556	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	21470	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	11982	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	23543m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13987m	0.40	ng/ul	-0.03
20) Perylene-d12	23.49	264	12187m	0.40	ng/ul	-0.05
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	5486	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	10736m	0.17	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.09	178	4222m	0.056	ng/ul	
15) Fluoranthene	19.12	202	7040m	0.084	ng/ul	
17) Pyrene	19.49	202	4682m	0.064	ng/ul	
18) Benzo(a)anthracene	21.25	228	1943m	0.031	ng/ul	
19) Chrysene	21.30	228	2894m	0.050	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	4458m	0.088	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	1941m	0.039	ng/ul	
23) Benzo(a)pyrene	23.39	252	1847	0.039	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.73	276	1924	0.038	ng/ul#	43
26) Benzo(g,h,i)perylene	26.41	276	1752	0.041	ng/ul#	55

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

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