

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006145.D
 Acq On : 06 Jun 2019 20:35
 Operator : JU/SJ
 Sample : K3201-02
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AB3

Manual Integrations
 APPROVED

mohammad
 6/11/2019 5:07:24 PM

Quant Time: Jun 07 01:18:08 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 : Fri Jun 07 00:25:28 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5929	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	22131	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	11632	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	24313m	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	21669m	0.40	ng/ul	-0.03
20) Perylene-d12	23.50	264	17827m	0.40	ng/ul	-0.06
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	9721	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	22333m	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.48	128	4860	0.080	ng/ul#	93
5) 2-Methylnaphthalene	12.10	142	2126	0.050	ng/ul	96
7) Acenaphthylene	14.02	152	8925	0.138	ng/ul	96
8) Acenaphthene	14.36	153	6114	0.136	ng/ul	98
9) Fluorene	15.36	166	7583	0.144	ng/ul	98
12) Phenanthrene	17.09	178	115541m	1.480	ng/ul	
13) Anthracene	17.19	178	28341m	0.380	ng/ul	
15) Fluoranthene	19.12	202	202724m	2.352	ng/ul	
17) Pyrene	19.49	202	171358	1.501	ng/ul	100
18) Benzo(a)anthracene	21.25	228	89814m	0.937	ng/ul	
19) Chrysene	21.30	228	94006	1.045	ng/ul	98
21) Benzo(b)fluoranthene	22.84	252	119512m	1.608	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	50697m	0.700	ng/ul	
23) Benzo(a)pyrene	23.41	252	80157	1.158	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.73	276	60896	0.813	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.74	278	14216	0.237	ng/ul	95
26) Benzo(a,h,i)perylene	26.42	276	46632	0.743	ng/ul	96

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

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