

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003092.D
 Acq On : 11 Oct 2018 07:39
 Operator : MJ/SJ
 Sample : J5295-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3Z48

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:45:05 PM

Quant Time: Oct 11 09:14:07 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN100918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 08:10:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2104	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	8054	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4125	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.01	188	7168m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	4890m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4949m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3225	0.26	ng/ul	-0.02
14) Fluoranthene-d10	19.05	212	5488m	0.27	ng/ul	-0.01
Target Compounds				Ovalue		
7) Acenaphthylene	13.98	152	1637	0.087	ng/ul#	55
12) Phenanthrene	17.06	178	6890m	0.345	ng/ul	
13) Anthracene	17.15	178	1696m	0.090	ng/ul	
15) Fluoranthene	19.09	202	22696m	0.968	ng/ul	
17) Pyrene	19.45	202	18864m	0.993	ng/ul	
18) Benzo(a)anthracene	21.21	228	10184	0.647	ng/ul#	79
19) Chrysene	21.26	228	10069	0.630	ng/ul#	78
21) Benzo(b)fluoranthene	22.79	252	12395m	0.695	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	4799m	0.260	ng/ul	
23) Benzo(a)pyrene	23.34	252	7944	0.497	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.67	276	6463	0.352	ng/ul#	78
26) Benzo(g,h,i)perylene	26.34	276	6631	0.428	ng/ul#	32

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

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