

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003065.D
 Acq On : 10 Oct 2018 15:20
 Operator : MJ/SJ
 Sample : J5295-08
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3Z42

Quant Time: Oct 11 07:07:26 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 07:04:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1452	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	5946	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3505	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7978	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	6207	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4845	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3324	0.36	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	9036	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	5499	0.360	ng/ul	99
5) 2-Methylnaphthalene	12.08	142	3668	0.349	ng/ul	97
7) Acenaphthylene	13.99	152	5738	0.358	ng/ul	99
9) Fluorene	15.33	166	6125	0.439	ng/ul	100
12) Phenanthrene	17.07	178	19996	0.899	ng/ul	99
15) Fluoranthene	19.09	202	14630	0.561	ng/ul	91
17) Pyrene	19.45	202	15515	0.644	ng/ul	99
18) Benzo(a)anthracene	21.21	228	7681	0.385	ng/ul	100
19) Chrysene	21.27	228	9877	0.487	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	6323	0.362	ng/ul	98
22) Benzo(k)fluoranthene	22.83	252	10366	0.573	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.66	276	8245	0.459	ng/ul#	96
25) Dibenzo(a,h)anthracene	25.66	278	7116	0.495	ng/ul	95
26) Benzo(g,h,i)perylene	26.34	276	10796	0.712	ng/ul	95

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

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