

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121918\
 Data File : BN004107.D
 Acq On : 19 Dec 2018 11:25
 Operator : JU/SJ
 Sample : J6271-13DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E28DL

Manual Integrations
 APPROVED

Sohil
 12/20/2018 2:02:47 PM

Quant Time: Dec 20 04:55:24 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 20 04:42:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1316	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	6670	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	4074	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	10558	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	11771	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	11517	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	481	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	1512	0.05	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	7544	0.411	ng/ul	98
12) Phenanthrene	17.25	178	1621m	0.047	ng/ul	
13) Anthracene	17.35	178	7409	0.260	ng/ul	98
15) Fluoranthene	19.27	202	30814	0.753	ng/ul	96
17) Pyrene	19.64	202	36194	0.857	ng/ul	99
18) Benzo(a)anthracene	21.38	228	22628	0.609	ng/ul	97
19) Chrysene	21.43	228	23601	0.555	ng/ul	97
21) Benzo(b)fluoranthene	23.02	252	45617m	0.958	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	13128m	0.283	ng/ul	
23) Benzo(a)pyrene	23.62	252	20492	0.497	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	26.08	276	18973	0.401	ng/ul#	94
25) Dibenzo(a,h)anthracene	26.08	278	4278m	0.112	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	14335	0.349	ng/ul	94

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

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Quant Time: Dec 20 04:55:24 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 20 04:42:11 2018
Response via : Initial Calibration

