

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121518\
 Data File : BN003975.D
 Acq On : 16 Dec 2018 01:09
 Operator : JU/SJ
 Sample : J6171-17DL 5X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y27DL

Manual Integrations
 APPROVED

Sohil
 12/17/2018 4:24:05 PM

Quant Time: Dec 16 02:24:17 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 : Sat Dec 15 23:04:58 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	1857	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9383	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5694	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	14507	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	18607	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	17476	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	705	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	2321	0.05	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.26	178	3031	0.064	ng/ul#	83
15) Fluoranthene	19.27	202	1204	0.021	ng/ul	77
17) Pyrene	19.64	202	5993m	0.090	ng/ul	
18) Benzo(a)anthracene	21.38	228	2520	0.043	ng/ul#	12
19) Chrysene	21.43	228	3468	0.052	ng/ul#	55
21) Benzo(b)fluoranthene	23.01	252	1615	0.022	ng/ul#	1
23) Benzo(a)pyrene	23.61	252	2398	0.038	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.06	276	1881	0.026	ng/ul#	98
26) Benzo(g,h,i)perylene	26.80	276	4240	0.068	ng/ul#	55

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

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Quant Time: Dec 16 02:24:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN11918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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