

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121018\
 Data File : BN003800.D
 Acq On : 11 Dec 2018 04:14
 Operator : JU/SJ
 Sample : J6077-18DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9Y03DL

Manual Integrations
 APPROVED

Sohil
 12/11/2018 6:12:34 PM

Quant Time: Dec 11 08:16:27 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 : Tue Dec 11 04:22:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.84	152	5544	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	21744	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	10228	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	19891	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14533	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	15974	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	1278	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	2175	0.04	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.26	178	9482	0.145	ng/ul	96
15) Fluoranthene	19.27	202	17268	0.224	ng/ul	96
17) Pyrene	19.64	202	14150	0.271	ng/ul	99
18) Benzo(a)anthracene	21.38	228	4462	0.097	ng/ul#	50
19) Chrysene	21.43	228	6259	0.119	ng/ul#	59
21) Benzo(b)fluoranthene	23.01	252	8966m	0.136	ng/ul	
22) Benzo(k)fluoranthene	23.05	252	3225m	0.050	ng/ul	
23) Benzo(a)pyrene	23.61	252	4585	0.080	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.06	276	4337	0.066	ng/ul#	97
26) Benzo(g,h,i)perylene	26.79	276	3866	0.068	ng/ul#	1

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

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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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