

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121818\
 Data File : BN004100.D
 Acq On : 19 Dec 2018 04:59
 Operator : JU/SJ
 Sample : J6271-09
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F9E24

Manual Integrations
 APPROVED

Sohil
 12/19/2018 4:16:59 PM

Quant Time: Dec 19 05:46:10 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 : Wed Dec 19 04:08:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1808	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	9002	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5482	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.21	188	14050m	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	14502m	0.40	ng/ul	0.00
20) Perylene-d12	23.72	264	13819m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	3285	0.24	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	10241m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.19	152	8727	0.354	ng/ul	99
12) Phenanthrene	17.25	178	3557m	0.077	ng/ul	
13) Anthracene	17.35	178	7868m	0.208	ng/ul	
15) Fluoranthene	19.27	202	38258m	0.703	ng/ul	
17) Pyrene	19.64	202	41737m	0.802	ng/ul	
18) Benzo(a)anthracene	21.38	228	28365m	0.619	ng/ul	
19) Chrysene	21.43	228	29465m	0.562	ng/ul	
21) Benzo(b)fluoranthene	23.02	252	56119m	0.982	ng/ul	
22) Benzo(k)fluoranthene	23.06	252	17860m	0.321	ng/ul	
23) Benzo(a)pyrene	23.62	252	23695m	0.479	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.08	276	22289m	0.393	ng/ul	
25) Dibenzo(a,h)anthracene	26.08	278	5520m	0.121	ng/ul	
26) Benzo(g,h,i)perylene	26.81	276	16360m	0.332	ng/ul	

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

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