

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003072.D
 Acq On : 10 Oct 2018 19:56
 Operator : MJ/SJ
 Sample : J5157-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD7

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:45:33 PM

Quant Time: Oct 11 07:46:21 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	1621	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6464	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3502	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	6608m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	4679m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4614m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3920	0.39	ng/ul	-0.01
14) Fluoranthene-d10	19.06	212	6496m	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	1016	0.061	ng/ul#	85
7) Acenaphthylene	13.98	152	1455	0.091	ng/ul#	91
9) Fluorene	15.32	166	346	0.025	ng/ul#	91
12) Phenanthrene	17.06	178	1570m	0.085	ng/ul	
13) Anthracene	17.15	178	2982m	0.171	ng/ul	
15) Fluoranthene	19.09	202	2703m	0.125	ng/ul	
17) Pyrene	19.45	202	10732m	0.591	ng/ul	
18) Benzo(a)anthracene	21.21	228	4366m	0.290	ng/ul	
19) Chrysene	21.26	228	11504	0.752	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	19873m	1.196	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	7180m	0.417	ng/ul	
23) Benzo(a)pyrene	23.34	252	12908	0.867	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.66	276	7750	0.453	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.65	278	2272	0.166	ng/ul#	83
26) Benzo(g,h,i)perylene	26.33	276	6353	0.440	ng/ul	99

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

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