

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003043.D
 Acq On : 10 Oct 2018 00:47
 Operator : MJ/SJ
 Sample : J5157-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD5

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:54 PM

Quant Time: Oct 10 02:33:34 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 : Wed Oct 10 01:34:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2008	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	8321	0.40	ng/ul	-0.03
6) Acenaphthene-d10	14.26	164	4898	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	10234m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7117m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6617	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	5392	0.42	ng/ul	-0.02
14) Fluoranthene-d10	19.05	212	11768m	0.40	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	1348	0.063	ng/ul#	90
7) Acenaphthylene	13.98	152	2005	0.089	ng/ul#	92
12) Phenanthrene	17.06	178	1350m	0.047	ng/ul	
13) Anthracene	17.15	178	2432m	0.090	ng/ul	
15) Fluoranthene	19.09	202	27148m	0.811	ng/ul	
17) Pyrene	19.45	202	32572m	1.178	ng/ul	
18) Benzo(a)anthracene	21.21	228	11779	0.514	ng/ul	98
19) Chrysene	21.26	228	13351	0.574	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	26408m	1.108	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	8547m	0.346	ng/ul	
23) Benzo(a)pyrene	23.34	252	16292	0.763	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.66	276	9157	0.373	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.65	278	2743	0.140	ng/ul#	82
26) Benzo(g,h,i)perylene	26.33	276	8631	0.417	ng/ul	99

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

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