

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101118\
 Data File : BN003080.D
 Acq On : 11 Oct 2018 00:37
 Operator : MJ/SJ
 Sample : J5157-08
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD0

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:45:53 PM

Quant Time: Oct 11 08:27:53 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 08:10:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1540	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6272	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3567	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.01	188	6871m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	4474m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4150m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	4371	0.45	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	8295m	0.42	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	1204	0.075	ng/ul#	87
7) Acenaphthylene	13.98	152	1889	0.116	ng/ul#	92
8) Acenaphthene	14.32	153	328	0.026	ng/ul#	94
9) Fluorene	15.32	166	436	0.031	ng/ul#	94
12) Phenanthrene	17.06	178	3202	0.167	ng/ul	97
13) Anthracene	17.15	178	3063m	0.169	ng/ul	
15) Fluoranthene	19.09	202	11252m	0.501	ng/ul	
17) Pyrene	19.44	202	24064m	1.385	ng/ul	
18) Benzo(a)anthracene	21.21	228	10847	0.753	ng/ul	98
19) Chrysene	21.26	228	12121	0.829	ng/ul	100
21) Benzo(b)fluoranthene	22.77	252	27576m	1.845	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	9815m	0.633	ng/ul	
23) Benzo(a)pyrene	23.33	252	16759	1.251	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.65	276	9881	0.642	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.65	278	2607	0.212	ng/ul	91
26) Benzo(a,h,i)perylene	26.33	276	7560	0.582	ng/ul	98

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

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