

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003036.D
 Acq On : 09 Oct 2018 20:41
 Operator : MJ/SJ
 Sample : J5157-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC8

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:47 PM

Quant Time: Oct 10 02:00:00 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	1491	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	6461	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4010	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9175m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	8670m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	7699	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	4306	0.43	ng/ul	-0.02
14) Fluoranthene-d10	19.06	212	11306m	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	1541	0.093	ng/ul#	90
5) 2-Methylnaphthalene	12.09	142	241	0.021	ng/ul	94
7) Acenaphthylene	13.98	152	1800	0.098	ng/ul#	93
8) Acenaphthene	14.32	153	344	0.025	ng/ul#	92
9) Fluorene	15.32	166	451	0.028	ng/ul#	93
12) Phenanthrene	17.06	178	1949m	0.076	ng/ul	
13) Anthracene	17.15	178	3007m	0.124	ng/ul	
15) Fluoranthene	19.09	202	3347m	0.112	ng/ul	
17) Pyrene	19.45	202	24534m	0.729	ng/ul	
18) Benzo(a)anthracene	21.21	228	6041m	0.217	ng/ul	
19) Chrysene	21.27	228	14709	0.519	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	17671m	0.637	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	7982m	0.278	ng/ul	
23) Benzo(a)pyrene	23.34	252	11069	0.445	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.66	276	5487	0.192	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.66	278	1880	0.082	ng/ul#	71
26) Benzo(g,h,i)perylene	26.34	276	5004	0.208	ng/ul	98

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

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