

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN101018\
 Data File : BN003041.D
 Acq On : 09 Oct 2018 23:37
 Operator : MJ/SJ
 Sample : J5157-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AD3

Manual Integrations
 APPROVED

Sohil
 10/10/2018 4:57:52 PM

Quant Time: Oct 10 02:25:33 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	1939	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	8093	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4683	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	9365m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	5885m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	6471	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	5244	0.42	ng/ul	-0.02
14) Fluoranthene-d10	19.06	212	10256m	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	2804	0.135	ng/ul	98
5) 2-Methylnaphthalene	12.08	142	315	0.022	ng/ul	99
7) Acenaphthylene	13.98	152	1632	0.076	ng/ul#	89
9) Fluorene	15.32	166	509	0.027	ng/ul#	92
12) Phenanthrene	17.06	178	2659m	0.102	ng/ul	
13) Anthracene	17.15	178	3477m	0.141	ng/ul	
15) Fluoranthene	19.09	202	6033m	0.197	ng/ul	
17) Pyrene	19.45	202	20156m	0.882	ng/ul	
18) Benzo(a)anthracene	21.21	228	10398	0.549	ng/ul	98
19) Chrysene	21.26	228	13589	0.707	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	24768m	1.063	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	8324m	0.345	ng/ul	
23) Benzo(a)pyrene	23.34	252	14877	0.712	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.66	276	7173	0.299	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.66	278	2134	0.111	ng/ul#	74
26) Benzo(g,h,i)perylene	26.33	276	5926	0.293	ng/ul	99

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

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