

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100918\
 Data File : BN003011.D
 Acq On : 09 Oct 2018 05:30
 Operator : MJ/SJ
 Sample : J5116-09
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLC58

Manual Integrations
 APPROVED

Sohil
 10/9/2018 5:38:23 PM

Quant Time: Oct 09 09:01:57 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 : Tue Oct 09 08:56:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1223	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	5247	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3165	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.02	188	7326m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	7461m	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	7609	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	2778	0.34	ng/ul	-0.01
14) Fluoranthene-d10	19.05	212	7454m	0.36	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	2157	0.160	ng/ul	95
5) 2-Methylnaphthalene	12.09	142	369	0.040	ng/ul	96
7) Acenaphthylene	13.98	152	497	0.034	ng/ul#	70
8) Acenaphthene	14.33	153	2325	0.212	ng/ul	99
9) Fluorene	15.32	166	1689	0.134	ng/ul#	97
12) Phenanthrene	17.06	178	37522	1.837	ng/ul	99
13) Anthracene	17.15	178	6607m	0.343	ng/ul	
15) Fluoranthene	19.09	202	89107m	3.719	ng/ul	
17) Pyrene	19.45	202	82233	2.838	ng/ul	97
18) Benzo(a)anthracene	21.21	228	43117	1.796	ng/ul	99
19) Chrysene	21.26	228	51296	2.104	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	69645m	2.542	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	23256m	0.819	ng/ul	
23) Benzo(a)pyrene	23.34	252	45026	1.833	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.65	276	27445	0.972	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.64	278	6425	0.284	ng/ul#	95
26) Benzo(g,h,i)perylene	26.32	276	24159	1.015	ng/ul	95

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

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