

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003114.D
 Acq On : 11 Oct 2018 21:47
 Operator : MJ/SJ
 Sample : J5368-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YZ7

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:38:22 PM

Quant Time: Oct 12 06:55:59 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 : Fri Oct 12 05:29:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1713	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7389	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4343	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.01	188	8837m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	5403m	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	4634m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3748	0.33	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	7384m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	952	0.050	ng/ul#	83
5) 2-Methylnaphthalene	12.08	142	634	0.049	ng/ul	98
7) Acenaphthylene	13.98	152	2051	0.103	ng/ul#	7
8) Acenaphthene	14.32	153	1100	0.073	ng/ul#	91
9) Fluorene	15.32	166	1576	0.091	ng/ul#	89
12) Phenanthrene	17.06	178	27127	1.101	ng/ul	99
13) Anthracene	17.15	178	7196m	0.309	ng/ul	
15) Fluoranthene	19.09	202	49652m	1.718	ng/ul	
17) Pyrene	19.45	202	43665	2.081	ng/ul	96
18) Benzo(a)anthracene	21.21	228	20531	1.181	ng/ul	97
19) Chrysene	21.26	228	20271	1.148	ng/ul	97
21) Benzo(b)fluoranthene	22.78	252	25401m	1.522	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	9645m	0.557	ng/ul	
23) Benzo(a)pyrene	23.34	252	18329	1.225	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.66	276	13203	0.768	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.65	278	3446	0.250	ng/ul#	66
26) Benzo(a,h,i)perylene	26.33	276	11673	0.805	ng/ul	98

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

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