

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003130.D
 Acq On : 12 Oct 2018 07:10
 Operator : MJ/SJ
 Sample : J5234-11
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AK9

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:38:42 PM

Quant Time: Oct 12 12:39:34 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 07:27:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1915	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.39	136	7626	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	4856	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	6699m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	4894m	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	4575m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	2873	0.24	ng/ul	-0.03
14) Fluoranthene-d10	19.05	212	5309m	0.28	ng/ul	-0.01
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	9959	0.509	ng/ul#	96
5) 2-Methylnaphthalene	12.06	142	36111	2.681	ng/ul	97
7) Acenaphthylene	13.98	152	4873	0.219	ng/ul#	46
8) Acenaphthene	14.32	153	9015	0.535	ng/ul	98
9) Fluorene	15.31	166	40989	2.123	ng/ul	98
12) Phenanthrene	17.05	178	70252	3.762	ng/ul	97
13) Anthracene	17.15	178	22097m	1.253	ng/ul	
15) Fluoranthene	19.08	202	45695m	2.086	ng/ul	
17) Pyrene	19.45	202	54504	2.867	ng/ul	98
18) Benzo(a)anthracene	21.21	228	18693m	1.187	ng/ul	
19) Chrysene	21.26	228	18676m	1.168	ng/ul	
21) Benzo(b)fluoranthene	22.78	252	17792m	1.080	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	6437m	0.377	ng/ul	
23) Benzo(a)pyrene	23.34	252	14500m	0.982	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.65	276	8958	0.528	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.64	278	2307m	0.170	ng/ul	
26) Benzo(a,h,i)perylene	26.32	276	8164	0.571	ng/ul	92

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

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