

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101218\
 Data File : BN003117.D
 Acq On : 11 Oct 2018 23:32
 Operator : MJ/SJ
 Sample : J5230-04DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YY0DL

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:38:26 PM

Quant Time: Oct 12 07:11:02 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	1711	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.40	136	7082	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3825	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	7729	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	5089	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5487	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	599	0.05	ng/ul	0.01
14) Fluoranthene-d10	19.05	212	1132	0.05	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	2025	0.111	ng/ul#	93
5) 2-Methylnaphthalene	12.09	142	1285	0.103	ng/ul	95
7) Acenaphthylene	13.98	152	15684	0.896	ng/ul	99
8) Acenaphthene	14.32	153	703	0.053	ng/ul#	93
9) Fluorene	15.32	166	1180	0.078	ng/ul#	96
12) Phenanthrene	17.06	178	27316	1.268	ng/ul	99
13) Anthracene	17.15	178	15645	0.769	ng/ul	99
15) Fluoranthene	19.08	202	123208	4.874	ng/ul	100
17) Pyrene	19.45	202	104145	5.269	ng/ul	97
18) Benzo(a)anthracene	21.21	228	49771	3.039	ng/ul	100
19) Chrysene	21.26	228	55232	3.321	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	80302m	4.064	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	31191m	1.523	ng/ul	
23) Benzo(a)pyrene	23.34	252	41666	2.353	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.64	276	34433	1.692	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.64	278	10006	0.614	ng/ul	100
26) Benzo(a,h,i)perylene	26.32	276	28406	1.655	ng/ul	94

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

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