

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN101218\
 Data File : BN003100.D
 Acq On : 11 Oct 2018 13:35
 Operator : MJ/SJ
 Sample : J5230-06
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3Z04

Quant Time: Oct 11 16:46:04 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 : Thu Oct 11 16:42:09 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	-0.02
2) Naphthalene-d8	10.39	136	6856	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.26	164	3726	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.11	188	345645	0.40	ng/ul	0.08
16) Chrysene-d12	21.27	240	77	0.40	ng/ul	0.03
20) Perylene-d12	23.44	264	4781	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3095	0.29	ng/ul	-0.02
14) Fluoranthene-d10	19.15	212	118	0.00	ng/ul	0.08
Target Compounds				Qvalue		
3) Naphthalene	10.44	128	14953	0.850	ng/ul	97
5) 2-Methylnaphthalene	12.07	142	12538	1.035	ng/ul	97
7) Acenaphthylene	13.98	152	29314	1.720	ng/ul	99
8) Acenaphthene	14.32	153	920	0.071	ng/ul#	92
9) Fluorene	15.32	166	1822	0.123	ng/ul#	81
12) Phenanthrene	17.15	178	25312	0.026	ng/ul	99
17) Pyrene	19.45	202	146418	489.587	ng/ul	95
18) Benzo(a)anthracene	21.26	228	103870	419.191	ng/ul	97
19) Chrysene	21.26	228	103552	411.558	ng/ul	100
21) Benzo(b)fluoranthene	22.77	252	210452	12.223	ng/ul	90
22) Benzo(k)fluoranthene	22.94	252	22061	1.236	ng/ul	98
23) Benzo(a)pyrene	23.34	252	73707	4.777	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.64	276	68817	3.881	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.64	278	19692	1.387	ng/ul	96
26) Benzo(g,h,i)perylene	26.31	276	57523	3.847	ng/ul	93

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

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