

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101420\
 Data File : BN012196.D
 Acq On : 14 Oct 2020 16:07
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV08

Quant Time: Oct 14 16:52:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 15:53:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1281	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	5017	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	2818	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5566	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	4772	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	4951	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	2919	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	5746	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	5648	0.390	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	3603	0.392	ng/ul	99
7) Acenaphthylene	13.98	152	4716	0.386	ng/ul	100
8) Acenaphthene	14.32	153	4006	0.388	ng/ul	98
9) Fluorene	15.32	166	4406	0.383	ng/ul	99
11) Pentachlorophenol	16.67	266	614	0.382	ng/ul	97
12) Phenanthrene	17.05	178	6889	0.388	ng/ul	99
13) Anthracene	17.15	178	5881	0.386	ng/ul	99
15) Fluoranthene	19.08	202	7922	0.382	ng/ul	99
17) Pyrene	19.45	202	8309	0.386	ng/ul	99
18) Benzo(a)anthracene	21.20	228	6672	0.377	ng/ul	99
19) Chrysene	21.25	228	8332	0.387	ng/ul	99
21) Benzo(b)fluoranthene	22.76	252	7770	0.389	ng/ul	99
22) Benzo(k)fluoranthene	22.80	252	9037	0.408	ng/ul	99
23) Benzo(a)pyrene	23.31	252	7626	0.392	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.60	276	9800	0.392	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.61	278	7779	0.389	ng/ul	98
26) Benzo(g,h,i)perylene	26.26	276	9000	0.393	ng/ul	100

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

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