

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112020\
 Data File : BN012661.D
 Acq On : 20 Nov 2020 15:51
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV07

Quant Time: Nov 20 16:39:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 15:24:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	768	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.20	136	2860	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.10	164	1560	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	3303	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	3361	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	3039	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	1738	0.42	ng/ul	-0.01
14) Fluoranthene-d10	18.89	212	3568	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.25	128	3518	0.409	ng/ul	99
5) 2-Methylnaphthalene	11.89	142	2236	0.407	ng/ul	99
7) Acenaphthylene	13.81	152	3059	0.402	ng/ul	98
8) Acenaphthene	14.16	153	2585	0.403	ng/ul	97
9) Fluorene	15.16	166	2895	0.411	ng/ul	97
11) Pentachlorophenol	16.51	266	424	0.404	ng/ul	96
12) Phenanthrene	16.89	178	4685	0.408	ng/ul	99
13) Anthracene	16.99	178	4061	0.406	ng/ul	99
15) Fluoranthene	18.93	202	5470	0.412	ng/ul	98
17) Pyrene	19.29	202	5604	0.373	ng/ul	97
18) Benzo(a)anthracene	21.06	228	4539	0.377	ng/ul	99
19) Chrysene	21.11	228	5708	0.379	ng/ul	99
21) Benzo(b)fluoranthene	22.57	252	5600	0.401	ng/ul	99
22) Benzo(k)fluoranthene	22.61	252	6220	0.386	ng/ul	98
23) Benzo(a)pyrene	23.10	252	5304	0.403	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.28	276	6966	0.405	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.29	278	5553	0.402	ng/ul	97
26) Benzo(g,h,i)perylene	25.91	276	6189	0.400	ng/ul	99

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

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