

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122818\
 Data File : BN004265.D
 Acq On : 28 Dec 2018 16:46
 Operator : JU/SJ
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV58

Quant Time: Dec 28 17:26:40 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN122818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 28 16:45:35 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	2067	0.40	ng/ul	0.00
2) Naphthalene-d8	10.62	136	9263	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	5217	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	11869	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	14046	0.40	ng/ul	0.00
20) Perylene-d12	23.74	264	12090	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	6038	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.25	212	14359	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.67	128	10825	0.415	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	7650	0.422	ng/ul	98
7) Acenaphthylene	14.20	152	10791	0.413	ng/ul	100
8) Acenaphthene	14.53	153	8640	0.413	ng/ul	99
9) Fluorene	15.52	166	9796	0.413	ng/ul	100
11) Pentachlorophenol	16.90	266	513	0.298	ng/ul	97
12) Phenanthrene	17.26	178	15391	0.399	ng/ul	99
13) Anthracene	17.36	178	13889	0.396	ng/ul	99
15) Fluoranthene	19.28	202	19190	0.413	ng/ul	99
17) Pyrene	19.64	202	19035	0.396	ng/ul	98
18) Benzo(a)anthracene	21.39	228	19390	0.406	ng/ul	100
19) Chrysene	21.44	228	21275	0.416	ng/ul	99
21) Benzo(b)fluoranthene	23.03	252	18819	0.410	ng/ul	100
22) Benzo(k)fluoranthene	23.07	252	20487	0.421	ng/ul	100
23) Benzo(a)pyrene	23.63	252	17696	0.406	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.11	276	18290	0.405	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.11	278	14985	0.404	ng/ul	99
26) Benzo(g,h,i)perylene	26.84	276	14995	0.394	ng/ul	99

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

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