

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031620\
 Data File : BN010229.D
 Acq On : 16 Mar 2020 15:47
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV007

Quant Time: Mar 16 16:18:00 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN031620.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 16 15:36:58 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1935	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	6213	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	3354	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	8492	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	7672	0.40	ng/ul	0.00
20) Perylene-d12	23.37	264	6967	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3970	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	10368	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	7039	0.357	ng/ul	100
5) 2-Methylnaphthalene	12.06	142	4855	0.367	ng/ul	100
7) Acenaphthylene	13.97	152	7039	0.395	ng/ul	99
8) Acenaphthene	14.32	153	5291	0.386	ng/ul	99
9) Fluorene	15.31	166	6268	0.385	ng/ul	98
11) Pentachlorophenol	16.65	266	588	0.355	ng/ul	99
12) Phenanthrene	17.04	178	11039	0.361	ng/ul	100
13) Anthracene	17.13	178	9428	0.351	ng/ul	99
15) Fluoranthene	19.06	202	13320	0.358	ng/ul	100
17) Pyrene	19.42	202	13312	0.365	ng/ul	99
18) Benzo(a)anthracene	21.18	228	11491	0.348	ng/ul	100
19) Chrysene	21.23	228	13745	0.365	ng/ul	99
21) Benzo(b)fluoranthene	22.73	252	12285	0.389	ng/ul	99
22) Benzo(k)fluoranthene	22.77	252	12601	0.379	ng/ul	100
23) Benzo(a)pyrene	23.28	252	10982	0.402	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.53	276	12740	0.372	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.54	278	10432	0.372	ng/ul	99
26) Benzo(g,h,i)perylene	26.17	276	11444	0.380	ng/ul	99

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

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