

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN081220\
 Data File : BN011419.D
 Acq On : 12 Aug 2020 14:19
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV07

Quant Time: Aug 12 14:51:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 12 14:05:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	904	0.40	ng/ul	0.00
2) Naphthalene-d8	10.19	136	2523	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.07	164	1260	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.84	188	2618	0.40	ng/ul	0.00
16) Chrysene-d12	21.06	240	2429	0.40	ng/ul	0.01
20) Perylene-d12	23.18	264	2376	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.79	152	1229	0.34	ng/ul	0.00
14) Fluoranthene-d10	18.88	212	2755	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.24	128	2598	0.353	ng/ul	96
5) 2-Methylnaphthalene	11.87	142	1705	0.367	ng/ul	100
7) Acenaphthylene	13.79	152	2326	0.367	ng/ul	97
8) Acenaphthene	14.14	153	1802	0.365	ng/ul	99
9) Fluorene	15.14	166	2275	0.391	ng/ul	99
11) Pentachlorophenol	16.51	266	236	0.380	ng/ul	94
12) Phenanthrene	16.88	178	3191	0.360	ng/ul	99
13) Anthracene	16.97	178	2687	0.369	ng/ul	97
15) Fluoranthene	18.91	202	4128	0.405	ng/ul	99
17) Pyrene	19.27	202	3950	0.416	ng/ul	99
18) Benzo(a)anthracene	21.04	228	2870	0.352	ng/ul	98
19) Chrysene	21.10	228	4128	0.376	ng/ul	100
21) Benzo(b)fluoranthene	22.55	252	3640	0.352	ng/ul	94
22) Benzo(k)fluoranthene	22.59	252	4156	0.364	ng/ul	94
23) Benzo(a)pyrene	23.08	252	3268	0.377	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.23	276	4432	0.378	ng/ul	99
25) Dibenzo(a,h)anthracene	25.25	278	3591	0.375	ng/ul	92
26) Benzo(g,h,i)perylene	25.85	276	3787	0.345	ng/ul	98

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

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