

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM040220\
 Data File : BM025892.D
 Acq On : 01 Apr 2020 17:58
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV83

Quant Time: Apr 01 18:30:26 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM040220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 17:54:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	2314	0.40	ng/ul	0.00
2) Naphthalene-d8	10.33	136	9444	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	5974	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	15493	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	16016	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	16475	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	6542	0.39	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	19318	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.38	128	10081	0.363	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	7441	0.383	ng/ul	99
7) Acenaphthylene	13.92	152	10549	0.413	ng/ul	98
8) Acenaphthene	14.27	153	8463	0.399	ng/ul	99
9) Fluorene	15.26	166	10505	0.400	ng/ul	100
11) Pentachlorophenol	16.61	266	1384	0.364	ng/ul	99
12) Phenanthrene	16.99	178	18256	0.374	ng/ul	99
13) Anthracene	17.08	178	15820	0.367	ng/ul	99
15) Fluoranthene	19.01	202	22075	0.366	ng/ul	99
17) Pyrene	19.37	202	22081	0.385	ng/ul	98
18) Benzo(a)anthracene	21.13	228	21230	0.371	ng/ul	100
19) Chrysene	21.18	228	23788	0.381	ng/ul	99
21) Benzo(b)fluoranthene	22.66	252	23955	0.395	ng/ul	100
22) Benzo(k)fluoranthene	22.70	252	24659	0.379	ng/ul	99
23) Benzo(a)pyrene	23.20	252	22023	0.406	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.41	276	27233	0.375	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.42	278	22810	0.377	ng/ul	99
26) Benzo(g,h,i)perylene	26.05	276	23029	0.375	ng/ul	100

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

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