

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030320\
 Data File : BM025212.D
 Acq On : 03 Mar 2020 14:36
 Operator : CG/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV99

Quant Time: Mar 03 15:20:58 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 03 14:30:08 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.68	152	2995	0.40	ng/ul	0.00
2) Naphthalene-d8	10.45	136	10870	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	7465	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	18317	0.40	ng/ul	0.00
16) Chrysene-d12	21.24	240	18708	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	17387	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.04	152	7278	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	20772	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.50	128	12072	0.369	ng/ul#	83
5) 2-Methylnaphthalene	12.12	142	8446	0.354	ng/ul	99
7) Acenaphthylene	14.02	152	12512	0.358	ng/ul	95
8) Acenaphthene	14.37	153	9858	0.364	ng/ul	100
9) Fluorene	15.36	166	12097	0.356	ng/ul	95
11) Pentachlorophenol	16.71	266	1521	0.317	ng/ul	97
12) Phenanthrene	17.09	178	21208	0.359	ng/ul	96
13) Anthracene	17.18	178	18359	0.350	ng/ul	95
15) Fluoranthene	19.11	202	26034	0.352	ng/ul	94
17) Pyrene	19.47	202	25559	0.360	ng/ul#	94
18) Benzo(a)anthracene	21.22	228	24278	0.353	ng/ul	98
19) Chrysene	21.27	228	27526	0.364	ng/ul	98
21) Benzo(b)fluoranthene	22.78	252	25697	0.369	ng/ul	92
22) Benzo(k)fluoranthene	22.82	252	26970	0.370	ng/ul#	95
23) Benzo(a)pyrene	23.34	252	21049	0.360	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.62	276	26224	0.360	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.63	278	21984	0.364	ng/ul#	94
26) Benzo(g,h,i)perylene	26.28	276	22384	0.362	ng/ul#	90

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

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