

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092320\
 Data File : BM027501.D
 Acq On : 23 Sep 2020 13:57
 Operator : JU/CG
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV083

Quant Time: Sep 23 14:28:25 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 23 13:53:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	1036	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	3023	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.45	164	2049	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.19	188	5130	0.40	ng/ul	0.00
16) Chrysene-d12	21.37	240	5114	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	5037	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	2143	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.22	212	6042	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	3568	0.403	ng/ul	98
5) 2-Methylnaphthalene	12.27	142	2715	0.422	ng/ul	99
7) Acenaphthylene	14.17	152	4381	0.419	ng/ul	100
8) Acenaphthene	14.52	153	3081	0.414	ng/ul	98
9) Fluorene	15.50	166	3821	0.415	ng/ul	96
11) Pentachlorophenol	16.84	266	648	0.464	ng/ul	99
12) Phenanthrene	17.24	178	6525	0.412	ng/ul	98
13) Anthracene	17.33	178	6372	0.411	ng/ul	99
15) Fluoranthene	19.25	202	8665	0.455	ng/ul	99
17) Pyrene	19.61	202	8593	0.368	ng/ul	99
18) Benzo(a)anthracene	21.36	228	8169	0.407	ng/ul	100
19) Chrysene	21.41	228	8598	0.414	ng/ul	99
21) Benzo(b)fluoranthene	22.97	252	8727	0.430	ng/ul	99
22) Benzo(k)fluoranthene	23.01	252	8813	0.418	ng/ul	99
23) Benzo(a)pyrene	23.55	252	7636	0.420	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.96	276	9735	0.416	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.97	278	7989	0.414	ng/ul	99
26) Benzo(g,h,i)perylene	26.65	276	8119	0.411	ng/ul	97

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

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