

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021821\
 Data File : BM028802.D
 Acq On : 18 Feb 2021 18:29
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.438

Quant Time: Feb 18 23:33:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 18 11:30:02 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	471	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1956	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	1044	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	2125	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	1967	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	1923	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1172	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2361	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	2041	0.419	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	1359	0.417	ng/ul	97
7) Acenaphthylene	14.26	152	1755	0.412	ng/ul	99
8) Acenaphthene	14.60	153	1386	0.410	ng/ul	99
9) Fluorene	15.59	166	1558	0.416	ng/ul	99
11) Pentachlorophenol	16.94	266	221	0.398	ng/ul	100
12) Phenanthrene	17.32	178	2433	0.410	ng/ul	99
13) Anthracene	17.41	178	2406	0.425	ng/ul	99
15) Fluoranthene	19.32	202	2810	0.416	ng/ul	98
17) Pyrene	19.68	202	2917	0.408	ng/ul	99
18) Benzo(a)anthracene	21.41	228	2703	0.403	ng/ul	99
19) Chrysene	21.46	228	2671	0.400	ng/ul	98
21) Benzo(b)fluoranthene	22.99	252	2744	0.417	ng/ul	85
22) Benzo(k)fluoranthene	23.03	252	2766	0.421	ng/ul#	84
23) Benzo(a)pyrene	23.56	252	2402	0.410	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	25.92	276	3161	0.414	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.93	278	2621	0.409	ng/ul	96
26) Benzo(g,h,i)perylene	26.61	276	2633	0.414	ng/ul	98

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

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