

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020821\
 Data File : BM028670.D
 Acq On : 08 Feb 2021 12:31
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDC0.483

Quant Time: Feb 08 13:26:37 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 : Mon Feb 08 13:26:08 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.71	136	1783	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	920	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	1804	0.40	ng/ul	0.00
16) Chrysene-d12	21.43	240	1702	0.40	ng/ul	0.00
20) Perylene-d12	23.67	264	1739	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1045	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	1941	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	1864	0.420	ng/ul	99
5) 2-Methylnaphthalene	12.38	142	1241	0.418	ng/ul	98
7) Acenaphthylene	14.27	152	1533	0.408	ng/ul	99
8) Acenaphthene	14.61	153	1240	0.417	ng/ul	99
9) Fluorene	15.60	166	1360	0.412	ng/ul	98
11) Pentachlorophenol	16.95	266	185	0.392	ng/ul	96
12) Phenanthrene	17.33	178	2111	0.419	ng/ul	99
13) Anthracene	17.42	178	1987	0.413	ng/ul	100
15) Fluoranthene	19.33	202	2371	0.413	ng/ul	97
17) Pyrene	19.69	202	2451	0.396	ng/ul	99
18) Benzo(a)anthracene	21.42	228	2362	0.407	ng/ul	98
19) Chrysene	21.47	228	2382	0.412	ng/ul	98
21) Benzo(b)fluoranthene	23.00	252	2501	0.420	ng/ul	94
22) Benzo(k)fluoranthene	23.04	252	2476	0.417	ng/ul	95
23) Benzo(a)pyrene	23.57	252	2211	0.417	ng/ul	93
24) Indeno(1,2,3-cd)pyrene	25.93	276	2881	0.417	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.94	278	2389	0.412	ng/ul	94
26) Benzo(g,h,i)perylene	26.63	276	2433	0.423	ng/ul	98

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

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