

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021621\
 Data File : BM028740.D
 Acq On : 16 Feb 2021 09:41
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.431

Quant Time: Feb 16 10:21:52 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 : Tue Feb 16 10:21:34 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	410	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	1630	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	860	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.28	188	1728	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	1548	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	1498	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	943	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	1777	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.75	128	1730	0.426	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	1135	0.418	ng/ul	98
7) Acenaphthylene	14.26	152	1345	0.383	ng/ul	99
8) Acenaphthene	14.60	153	1159	0.417	ng/ul	98
9) Fluorene	15.59	166	1293	0.419	ng/ul	98
11) Pentachlorophenol	16.94	266	129	0.286	ng/ul	96
12) Phenanthrene	17.32	178	2013	0.417	ng/ul	100
13) Anthracene	17.41	178	1794	0.390	ng/ul	100
15) Fluoranthene	19.32	202	2186	0.398	ng/ul	98
17) Pyrene	19.68	202	2217	0.394	ng/ul	99
18) Benzo(a)anthracene	21.41	228	2008	0.380	ng/ul	98
19) Chrysene	21.46	228	2155	0.410	ng/ul	98
21) Benzo(b)fluoranthene	22.98	252	2255	0.440	ng/ul	88
22) Benzo(k)fluoranthene	23.03	252	2192	0.428	ng/ul#	90
23) Benzo(a)pyrene	23.56	252	1888	0.413	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.91	276	2524	0.424	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.92	278	2075	0.416	ng/ul#	91
26) Benzo(g,h,i)perylene	26.60	276	2175	0.439	ng/ul	96

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

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