

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021821\
 Data File : BM028788.D
 Acq On : 18 Feb 2021 09:59
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.437

Quant Time: Feb 18 11:30:21 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	458	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1936	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	1029	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	2066	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	1914	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	1910	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1157	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2256	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	2001	0.415	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	1343	0.416	ng/ul	99
7) Acenaphthylene	14.26	152	1727	0.411	ng/ul	99
8) Acenaphthene	14.60	153	1390	0.418	ng/ul	98
9) Fluorene	15.58	166	1581	0.428	ng/ul	98
11) Pentachlorophenol	16.94	266	223	0.413	ng/ul	97
12) Phenanthrene	17.31	178	2408	0.417	ng/ul	98
13) Anthracene	17.41	178	2354	0.427	ng/ul	99
15) Fluoranthene	19.32	202	2739	0.417	ng/ul	99
17) Pyrene	19.68	202	2832	0.407	ng/ul	98
18) Benzo(a)anthracene	21.41	228	2633	0.403	ng/ul	98
19) Chrysene	21.46	228	2637	0.406	ng/ul	98
21) Benzo(b)fluoranthene	22.98	252	2718	0.416	ng/ul	87
22) Benzo(k)fluoranthene	23.03	252	2729	0.418	ng/ul#	88
23) Benzo(a)pyrene	23.56	252	2397	0.412	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	25.91	276	3140	0.414	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.92	278	2580	0.406	ng/ul#	91
26) Benzo(g,h,i)perylene	26.61	276	2613	0.414	ng/ul	99

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

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