

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021821\
 Data File : BM028811.D
 Acq On : 19 Feb 2021 01:07
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.439

Quant Time: Feb 19 10:09:47 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 : Fri Feb 19 10:07:28 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	520	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	2175	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	1120	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	2221	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	2072	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	2091	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	1276	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2436	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	2257	0.417	ng/ul	100
5) 2-Methylnaphthalene	12.36	142	1496	0.413	ng/ul	98
7) Acenaphthylene	14.26	152	1856	0.406	ng/ul	100
8) Acenaphthene	14.60	153	1493	0.412	ng/ul	99
9) Fluorene	15.58	166	1667	0.414	ng/ul	97
11) Pentachlorophenol	16.94	266	222	0.383	ng/ul	96
12) Phenanthrene	17.31	178	2579	0.415	ng/ul	99
13) Anthracene	17.41	178	2512	0.424	ng/ul	100
15) Fluoranthene	19.32	202	2952	0.418	ng/ul	97
17) Pyrene	19.68	202	3036	0.403	ng/ul	98
18) Benzo(a)anthracene	21.40	228	2905	0.411	ng/ul	99
19) Chrysene	21.45	228	2895	0.412	ng/ul	99
21) Benzo(b)fluoranthene	22.98	252	2960	0.414	ng/ul	91
22) Benzo(k)fluoranthene	23.02	252	3073	0.430	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	2631	0.413	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.90	276	3460	0.417	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	2862	0.411	ng/ul	96
26) Benzo(g,h,i)perylene	26.60	276	2905	0.420	ng/ul	98

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

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