

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM121120\
 Data File : BM028103.D
 Acq On : 11 Dec 2020 17:18
 Operator : JU/CG
 Sample : SSTD0.104
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.1041

Quant Time: Dec 12 02:57:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SFAM-EPA-SIM-BM121120.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 12 02:45:43 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	3092	0.40	ng/ul	0.00
4) Naphthalene-d8	11.05	136	11636	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.85	164	5959	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.56	188	11684	0.40	ng/ul	0.00
17) Chrysene-d12	21.67	240	9260	0.40	ng/ul	0.00
23) Perylene-d12	24.04	264	8223	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/ul	
6) 2-Methylnaphthalene-d10	12.62	152	1739	0.09	ng/ul	0.00
18) Fluoranthene-d10	19.55	212	2836	0.10	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	11.10	128	3331	0.101	ng/ul	96
7) 2-Methylnaphthalene	12.69	142	2109	0.093	ng/ul	99
8) 1-Methylnaphthalene	12.91	142	2009	0.094	ng/ul	98
10) Acenaphthylene	14.57	152	2383	0.090	ng/ul#	95
11) Acenaphthene	14.91	153	2124	0.105	ng/ul	98
12) Fluorene	15.88	166	2343	0.098	ng/ul	98
15) Phenanthrene	17.60	178	3649	0.107	ng/ul	96
16) Anthracene	17.69	178	3130	0.094	ng/ul	95
19) Fluoranthene	19.58	202	3745	0.103	ng/ul#	91
20) Pyrene	19.94	202	3823	0.106	ng/ul#	90
21) Benzo(a)anthracene	21.66	228	3155	0.093	ng/ul	97
22) Chrysene	21.71	228	3416	0.098	ng/ul	98
24) Benzo(b)fluoranthene	23.32	252	3356	0.102	ng/ul#	62
25) Benzo(k)fluoranthene	23.37	252	3067	0.093	ng/ul#	62
26) Benzo(a)pyrene	23.94	252	2668	0.089	ng/ul#	52
27) Indeno(1,2,3-cd)pyrene	26.47	276	3469	0.099	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.48	278	2905	0.099	ng/ul#	76
29) Benzo(a,h,i)perylene	27.23	276	3002	0.101	ng/ul#	92

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

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