

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM022420\
 Data File : BM025032.D
 Acq On : 24 Feb 2020 21:23
 Operator : CG/JU
 Sample : SSTD1.681
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.681

Quant Time: Feb 25 00:51:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM022420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 25 00:46:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	1807	0.40	ng/ul	0.00
2) Naphthalene-d8	10.47	136	7299	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.33	164	4956	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	11813	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	13043	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	13769	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.07	152	21583	1.80	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	64542	1.99	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.52	128	34940	1.680	ng/ul#	85
5) 2-Methylnaphthalene	12.14	142	25602	1.796	ng/ul	100
7) Acenaphthylene	14.04	152	45072	1.965	ng/ul	94
8) Acenaphthene	14.39	153	30113	1.568	ng/ul	98
9) Fluorene	15.38	166	37256	1.693	ng/ul	96
11) Pentachlorophenol	16.72	266	6008	2.507	ng/ul	99
12) Phenanthrene	17.11	178	62132	1.651	ng/ul	96
13) Anthracene	17.20	178	63442	1.965	ng/ul	96
15) Fluoranthene	19.13	202	81150	1.998	ng/ul	97
17) Pyrene	19.49	202	82434	1.394	ng/ul	97
18) Benzo(a)anthracene	21.24	228	83122	1.665	ng/ul	98
19) Chrysene	21.29	228	81765	1.557	ng/ul	98
21) Benzo(b)fluoranthene	22.80	252	85764	1.434	ng/ul	92
22) Benzo(k)fluoranthene	22.84	252	84739	1.388	ng/ul#	93
23) Benzo(a)pyrene	23.36	252	79218	1.513	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.65	276	99271	1.555	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.67	278	83162	1.612	ng/ul#	94
26) Benzo(g,h,i)perylene	26.32	276	83410	1.640	ng/ul#	92

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

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