

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112020\
 Data File : BM027853.D
 Acq On : 20 Nov 2020 15:19
 Operator : JU/CG
 Sample : SSTD1.677
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.677

Quant Time: Nov 20 17:15:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 20 17:13:49 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.34	152	844	0.40	ng/ul	0.00
2) Naphthalene-d8	11.17	136	3298	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	2032	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.67	188	4306	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	3168	0.40	ng/ul	0.00
20) Perylene-d12	24.19	264	2415	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	9167	1.65	ng/ul	0.00
14) Fluoranthene-d10	19.65	212	19612	1.70	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	15620	1.580	ng/ul	97
5) 2-Methylnaphthalene	12.81	142	10753	1.599	ng/ul	99
7) Acenaphthylene	14.68	152	16204	1.634	ng/ul	97
8) Acenaphthene	15.01	153	11893	1.601	ng/ul	99
9) Fluorene	15.98	166	13984	1.622	ng/ul	99
11) Pentachlorophenol	17.31	266	2436	1.661	ng/ul	97
12) Phenanthrene	17.70	178	21899	1.610	ng/ul	98
13) Anthracene	17.79	178	21427	1.633	ng/ul	99
15) Fluoranthene	19.68	202	25073	1.662	ng/ul	99
17) Pyrene	20.04	202	24845	1.544	ng/ul	98
18) Benzo(a)anthracene	21.75	228	20403	1.596	ng/ul	99
19) Chrysene	21.81	228	19858	1.595	ng/ul	99
21) Benzo(b)fluoranthene	23.45	252	16494	1.625	ng/ul	92
22) Benzo(k)fluoranthene	23.50	252	16178	1.628	ng/ul#	93
23) Benzo(a)pyrene	24.08	252	13996	1.577	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	26.69	276	14715	1.532	ng/ul#	99
25) Dibenzo(a,h)anthracene	26.70	278	12298	1.537	ng/ul	92
26) Benzo(g,h,i)perylene	27.46	276	12287	1.524	ng/ul	95

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

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