

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030920\
 Data File : BM025413.D
 Acq On : 09 Mar 2020 12:22
 Operator : CG/JU
 Sample : SSTD1.677
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD1.677

Quant Time: Mar 09 13:26:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM030920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 09 12:03:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	2007	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	7625	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	5034	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	10344	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	8637	0.40	ng/ul	0.00
20) Perylene-d12	23.41	264	7705	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	23240	1.63	ng/ul	0.00
14) Fluoranthene-d10	19.06	212	49406	1.47	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.48	128	37837	1.649	ng/ul	99
5) 2-Methylnaphthalene	12.10	142	27190	1.627	ng/ul	99
7) Acenaphthylene	14.01	152	34827	1.477	ng/ul	99
8) Acenaphthene	14.35	153	30585	1.673	ng/ul	100
9) Fluorene	15.34	166	36999	1.614	ng/ul	99
11) Pentachlorophenol	16.69	266	4513	1.667	ng/ul	99
12) Phenanthrene	17.07	178	56297	1.688	ng/ul	100
13) Anthracene	17.16	178	51132	1.727	ng/ul	99
15) Fluoranthene	19.09	202	62661	1.501	ng/ul	100
17) Pyrene	19.45	202	59971	1.831	ng/ul	99
18) Benzo(a)anthracene	21.20	228	52962	1.668	ng/ul	100
19) Chrysene	21.26	228	57489	1.648	ng/ul	98
21) Benzo(b)fluoranthene	22.76	252	52838	1.714	ng/ul	94
22) Benzo(k)fluoranthene	22.80	252	55579	1.719	ng/ul#	95
23) Benzo(a)pyrene	23.31	252	45147	1.744	ng/ul#	92
24) Indeno(1,2,3-cd)pyrene	25.58	276	62404	1.934	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.59	278	53205	1.990	ng/ul	96
26) Benzo(g,h,i)perylene	26.24	276	54026	1.972	ng/ul	98

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

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