

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM041520\
 Data File : BM025957.D
 Acq On : 15 Apr 2020 09:27
 Operator : MA/SJ
 Sample : SSTD0.281
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.281

Quant Time: Apr 15 10:44:56 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM041520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 15 10:41:32 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1052	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	3807	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	2290	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	5162	0.40	ng/ul	0.00
16) Chrysene-d12	21.14	240	3986	0.40	ng/ul	0.00
20) Perylene-d12	23.29	264	3358	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	1319	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	3056	0.19	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	2537	0.227	ng/ul#	94
5) 2-Methylnaphthalene	12.00	142	1583	0.202	ng/ul	92
7) Acenaphthylene	13.92	152	2220	0.227	ng/ul	95
8) Acenaphthene	14.26	153	1914	0.236	ng/ul	95
9) Fluorene	15.26	166	2191	0.217	ng/ul	96
11) Pentachlorophenol	16.62	266	197	0.155	ng/ul	91
12) Phenanthrene	16.99	178	3696	0.227	ng/ul	98
13) Anthracene	17.08	178	3016	0.210	ng/ul	96
15) Fluoranthene	19.01	202	4150	0.207	ng/ul	96
17) Pyrene	19.37	202	4256	0.298	ng/ul	98
18) Benzo(a)anthracene	21.13	228	3071	0.216	ng/ul	97
19) Chrysene	21.18	228	4007	0.258	ng/ul	99
21) Benzo(b)fluoranthene	22.65	252	3157	0.255	ng/ul	92
22) Benzo(k)fluoranthene	22.69	252	3541	0.267	ng/ul#	92
23) Benzo(a)pyrene	23.19	252	2685	0.243	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	25.42	276	2869	0.194	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.44	278	2218	0.180	ng/ul#	88
26) Benzo(g,h,i)perylene	26.05	276	2702	0.216	ng/ul	93

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

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