

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM092320\
 Data File : BM027500.D
 Acq On : 23 Sep 2020 13:10
 Operator : JU/CG
 Sample : SSTD3.281
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD3.281

Quant Time: Sep 23 13:50:51 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM092320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 23 11:54:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.82	152	1309	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	3954	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.45	164	2744	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.20	188	6583	0.40	ng/ul	0.00
16) Chrysene-d12	21.37	240	6527	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	6211	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	22883	3.43	ng/ul	0.00
14) Fluoranthene-d10	19.22	212	64391	3.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	35654	3.011	ng/ul	97
5) 2-Methylnaphthalene	12.27	142	26139	3.223	ng/ul	99
7) Acenaphthylene	14.17	152	44118	3.703	ng/ul	97
8) Acenaphthene	14.51	153	30978	2.993	ng/ul	97
9) Fluorene	15.50	166	38885	3.134	ng/ul	99
11) Pentachlorophenol	16.84	266	6539	3.012	ng/ul	99
12) Phenanthrene	17.23	178	64604	2.902	ng/ul	99
13) Anthracene	17.33	178	65463	3.239	ng/ul	98
15) Fluoranthene	19.25	202	85827	3.208	ng/ul	98
17) Pyrene	19.61	202	85669	3.070	ng/ul	98
18) Benzo(a)anthracene	21.36	228	81347	3.137	ng/ul	100
19) Chrysene	21.41	228	82120	2.926	ng/ul	99
21) Benzo(b)fluoranthene	22.97	252	82356	3.092	ng/ul	95
22) Benzo(k)fluoranthene	23.01	252	86382	3.118	ng/ul#	96
23) Benzo(a)pyrene	23.55	252	75361	3.124	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	25.95	276	99251	2.983	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.97	278	81885	3.024	ng/ul#	95
26) Benzo(g,h,i)perylene	26.65	276	83047	2.892	ng/ul	98

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

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