

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092320\
 Data File : BM027497.D
 Acq On : 23 Sep 2020 11:21
 Operator : JU/CG
 Sample : SSTD0.478
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.478

Quant Time: Sep 23 11:55:25 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	1383	0.40	ng/ul	0.00
2) Naphthalene-d8	10.61	136	4121	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.45	164	2788	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.20	188	6649	0.40	ng/ul	0.00
16) Chrysene-d12	21.37	240	6107	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	6292	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.20	152	3029	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.22	212	7937	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.66	128	4854	0.393	ng/ul	100
5) 2-Methylnaphthalene	12.27	142	3501	0.414	ng/ul	100
7) Acenaphthylene	14.17	152	5768	0.476	ng/ul	100
8) Acenaphthene	14.51	153	4110	0.391	ng/ul	100
9) Fluorene	15.50	166	5031	0.399	ng/ul	100
11) Pentachlorophenol	16.84	266	731	0.333	ng/ul	100
12) Phenanthrene	17.24	178	8309	0.370	ng/ul	100
13) Anthracene	17.33	178	8128	0.398	ng/ul	100
15) Fluoranthene	19.25	202	10621	0.393	ng/ul	100
17) Pyrene	19.61	202	10571	0.405	ng/ul	100
18) Benzo(a)anthracene	21.36	228	9731	0.401	ng/ul	100
19) Chrysene	21.41	228	10062	0.383	ng/ul	100
21) Benzo(b)fluoranthene	22.97	252	10119	0.375	ng/ul	100
22) Benzo(k)fluoranthene	23.01	252	10602	0.378	ng/ul	100
23) Benzo(a)pyrene	23.55	252	9016	0.369	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.95	276	11988	0.356	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.97	278	9966	0.363	ng/ul	100
26) Benzo(g,h,i)perylene	26.65	276	10024	0.345	ng/ul	100

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

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