

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081820\
 Data File : BM027117.D
 Acq On : 18 Aug 2020 08:42
 Operator : JU/CG
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.425

Quant Time: Aug 18 11:17:49 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM081720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 17 14:20:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	1443	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	4419	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.26	164	2677	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	5964	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	5686	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	5585	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3172	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	7080	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	5718	0.432	ng/ul	100
5) 2-Methylnaphthalene	12.07	142	3903	0.431	ng/ul	95
7) Acenaphthylene	13.98	152	4822	0.415	ng/ul	99
8) Acenaphthene	14.33	153	4398	0.435	ng/ul	98
9) Fluorene	15.32	166	5219	0.431	ng/ul	98
11) Pentachlorophenol	16.67	266	712	0.362	ng/ul	94
12) Phenanthrene	17.05	178	8717	0.432	ng/ul	99
13) Anthracene	17.14	178	7658	0.418	ng/ul	99
15) Fluoranthene	19.07	202	10271	0.424	ng/ul	99
17) Pyrene	19.43	202	10327	0.425	ng/ul	99
18) Benzo(a)anthracene	21.19	228	9273	0.411	ng/ul	99
19) Chrysene	21.24	228	10849	0.444	ng/ul	99
21) Benzo(b)fluoranthene	22.74	252	10325	0.431	ng/ul	99
22) Benzo(k)fluoranthene	22.79	252	11035	0.443	ng/ul	100
23) Benzo(a)pyrene	23.30	252	9201	0.424	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.58	276	12728	0.425	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.60	278	10451	0.429	ng/ul	98
26) Benzo(g,h,i)perylene	26.25	276	11076	0.429	ng/ul	98

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

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