

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM020521\
 Data File : BM028663.D
 Acq On : 05 Feb 2021 14:26
 Operator : CG/JU
 Sample : SST0.476
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.476

Quant Time: Feb 05 15:58:43 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM020521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 05 15:47:53 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	475	0.40	ng/ul	0.00
2) Naphthalene-d8	10.72	136	1944	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	994	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.29	188	1928	0.40	ng/ul	0.00
16) Chrysene-d12	21.44	240	1685	0.40	ng/ul	0.00
20) Perylene-d12	23.68	264	1772	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1146	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.31	212	2043	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.76	128	2020	0.347	ng/ul	100
5) 2-Methylnaphthalene	12.38	142	1336	0.337	ng/ul	100
7) Acenaphthylene	14.28	152	1696	0.350	ng/ul	100
8) Acenaphthene	14.62	153	1327	0.365	ng/ul	100
9) Fluorene	15.60	166	1466	0.348	ng/ul	100
11) Pentachlorophenol	16.95	266	205	0.312	ng/ul	99
12) Phenanthrene	17.34	178	2227	0.366	ng/ul	100
13) Anthracene	17.43	178	2140	0.364	ng/ul	100
15) Fluoranthene	19.34	202	2470	0.366	ng/ul	100
17) Pyrene	19.70	202	2568	0.300	ng/ul	100
18) Benzo(a)anthracene	21.43	228	2373	0.349	ng/ul	100
19) Chrysene	21.48	228	2365	0.357	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	2482	0.333	ng/ul	100
22) Benzo(k)fluoranthene	23.05	252	2463	0.338	ng/ul	100
23) Benzo(a)pyrene	23.59	252	2185	0.336	ng/ul	100
24) Indeno(1,2,3-cd)pyrene	25.96	276	2900	0.411	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.97	278	2395	0.408	ng/ul	100
26) Benzo(g,h,i)perylene	26.66	276	2433	0.411	ng/ul	100

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

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