

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021621\
 Data File : BM028761.D
 Acq On : 16 Feb 2021 23:15
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.433

Quant Time: Feb 17 06:00:41 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 : Tue Feb 16 17:19:06 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	466	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	1850	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	976	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	1985	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	1802	0.40	ng/ul	0.00
20) Perylene-d12	23.65	264	1800	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1097	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.29	212	2078	0.41	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	1973	0.428	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	1280	0.415	ng/ul	99
7) Acenaphthylene	14.26	152	1540	0.387	ng/ul	98
8) Acenaphthene	14.60	153	1310	0.415	ng/ul	99
9) Fluorene	15.59	166	1483	0.423	ng/ul	99
11) Pentachlorophenol	16.94	266	178	0.343	ng/ul	98
12) Phenanthrene	17.31	178	2322	0.419	ng/ul	99
13) Anthracene	17.41	178	2091	0.395	ng/ul	99
15) Fluoranthene	19.32	202	2559	0.405	ng/ul	96
17) Pyrene	19.68	202	2598	0.397	ng/ul	99
18) Benzo(a)anthracene	21.40	228	2387	0.388	ng/ul	97
19) Chrysene	21.46	228	2502	0.409	ng/ul	99
21) Benzo(b)fluoranthene	22.98	252	2667	0.433	ng/ul	88
22) Benzo(k)fluoranthene	23.02	252	2594	0.422	ng/ul#	89
23) Benzo(a)pyrene	23.55	252	2245	0.409	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.90	276	3040	0.425	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.91	278	2535	0.423	ng/ul#	90
26) Benzo(g,h,i)perylene	26.60	276	2608	0.439	ng/ul	95

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

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