

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072221\
 Data File : BM031082.D
 Acq On : 23 Jul 2021 01:46
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4060

Quant Time: Jul 23 02:37:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 11:33:47 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.635	152	1096	0.400	ng/ul	0.00
4) Naphthalene-d8	10.396	136	4275	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.261	164	2835	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.009	188	6222	0.400	ng/ul	0.00
17) Chrysene-d12	21.193	240	5789	0.400	ng/ul	0.00
23) Perylene-d12	23.365	264	5373	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	416	0.361	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	2820	0.376	ng/ul	0.00
18) Fluoranthene-d10	19.039	212	6900	0.391	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.248	88	418	0.371	ng/ul	98
5) Naphthalene	10.446	128	4750	0.372	ng/ul	100
7) 2-Methylnaphthalene	12.065	142	3432	0.373	ng/ul	100
8) 1-Methylnaphthalene	12.285	142	3409	0.380	ng/ul	100
10) Acenaphthylene	13.981	152	4382	0.378	ng/ul	99
11) Acenaphthene	14.326	153	3820	0.373	ng/ul	97
12) Fluorene	15.315	166	4447	0.369	ng/ul	99
14) Pentachlorophenol	16.659	266	793	0.368	ng/ul	99
15) Phenanthrene	17.047	178	7421	0.376	ng/ul	98
16) Anthracene	17.141	178	6813	0.368	ng/ul	98
19) Fluoranthene	19.069	202	8847	0.380	ng/ul	99
20) Pyrene	19.432	202	9017	0.379	ng/ul	99
21) Benzo(a)anthracene	21.178	228	8694	0.376	ng/ul	99
22) Chrysene	21.229	228	9177	0.367	ng/ul	99
24) Benzo(b)fluoranthene	22.718	252	9508	0.376	ng/ul	99
25) Benzo(k)fluoranthene	22.759	252	9462	0.369	ng/ul	98
26) Benzo(a)pyrene	23.270	252	8000	0.366	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.519	276	10344	0.362	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.538	278	8205	0.357	ng/ul	98
29) Benzo(g,h,i)perylene	26.180	276	8724	0.359	ng/ul	99

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

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