

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038571.D
 Acq On : 27 Jan 2023 05:51
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4067

Quant Time: Jan 27 06:25:39 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM012623.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 27 00:18:15 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.967	152	1363	0.400	ng/ul	0.00
4) Naphthalene-d8	10.774	136	3624	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.597	164	2139	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.341	188	3946	0.400	ng/ul	0.00
17) Chrysene-d12	21.512	240	3742	0.400	ng/ul	0.00
23) Perylene-d12	23.924	264	4760	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.343	96	682	0.381	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.357	152	1923	0.342	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	3523	0.330	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	683	0.376	ng/ul#	23
5) Naphthalene	10.823	128	3323	0.282	ng/ul	99
7) 2-Methylnaphthalene	12.434	142	2166	0.348	ng/ul	97
8) 1-Methylnaphthalene	12.649	142	2164	0.337	ng/ul	98
10) Acenaphthylene	14.319	152	3305	0.350	ng/ul	96
11) Acenaphthene	14.657	153	2422	0.337	ng/ul	99
12) Fluorene	15.647	166	2719	0.303	ng/ul	98
14) Pentachlorophenol	16.999	266	554	0.345	ng/ul	94
15) Phenanthrene	17.384	178	3817	0.299	ng/ul	99
16) Anthracene	17.476	178	3644	0.335	ng/ul	97
19) Fluoranthene	19.394	202	4486	0.336	ng/ul	99
20) Pyrene	19.757	202	4746	0.331	ng/ul	98
21) Benzo(a)anthracene	21.498	228	4413	0.332	ng/ul	99
22) Chrysene	21.550	228	4544	0.340	ng/ul	98
24) Benzo(b)fluoranthene	23.187	252	5519	0.308	ng/ul	96
25) Benzo(k)fluoranthene	23.237	252	5682	0.327	ng/ul	97
26) Benzo(a)pyrene	23.819	252	5380	0.335	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.431	276	8288	0.366	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.444	278	6573	0.365	ng/ul	97
29) Benzo(g,h,i)perylene	27.212	276	7356	0.374	ng/ul	97

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

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