

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012623\
 Data File : BM038603.D
 Acq On : 28 Jan 2023 02:05
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4069

Quant Time: Jan 28 02:37:30 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	7.962	152	1445	0.400	ng/ul	0.00
4) Naphthalene-d8	10.768	136	4134	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.592	164	2703	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.337	188	5739	0.400	ng/ul	0.00
17) Chrysene-d12	21.510	240	5624	0.400	ng/ul	0.00
23) Perylene-d12	23.915	264	6333	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.338	96	692	0.365	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.352	152	2353	0.367	ng/ul	-0.01
18) Fluoranthene-d10	19.357	212	5922	0.369	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.376	88	698	0.362	ng/ul#	21
5) Naphthalene	10.817	128	3989	0.297	ng/ul	100
7) 2-Methylnaphthalene	12.429	142	2585	0.364	ng/ul	98
8) 1-Methylnaphthalene	12.643	142	2697	0.368	ng/ul	100
10) Acenaphthylene	14.319	152	4344	0.364	ng/ul	99
11) Acenaphthene	14.657	153	3200	0.352	ng/ul	97
12) Fluorene	15.642	166	3758	0.331	ng/ul	100
14) Pentachlorophenol	16.995	266	806	0.345	ng/ul	98
15) Phenanthrene	17.380	178	5824	0.314	ng/ul	99
16) Anthracene	17.468	178	5553	0.351	ng/ul	98
19) Fluoranthene	19.390	202	7536	0.376	ng/ul	99
20) Pyrene	19.752	202	7940	0.368	ng/ul	98
21) Benzo(a)anthracene	21.492	228	7119	0.356	ng/ul	98
22) Chrysene	21.545	228	7171	0.357	ng/ul	99
24) Benzo(b)fluoranthene	23.179	252	8150	0.342	ng/ul	98
25) Benzo(k)fluoranthene	23.228	252	8055	0.348	ng/ul	99
26) Benzo(a)pyrene	23.813	252	7407	0.346	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.425	276	10967	0.364	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.441	278	8715	0.364	ng/ul	99
29) Benzo(g,h,i)perylene	27.196	276	9485	0.363	ng/ul	98

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

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