

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011223\
 Data File : BM038437.D
 Acq On : 12 Jan 2023 13:01
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4064

Quant Time: Jan 12 14:14:34 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM123022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 12 01:15:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.996	152	1264	0.400	ng/u1	0.00
4) Naphthalene-d8	10.806	136	3724	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.624	164	2114	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.371	188	4444	0.400	ng/u1	0.00
17) Chrysene-d12	21.544	240	4501	0.400	ng/u1	0.00
23) Perylene-d12	23.974	264	4799	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	724	0.384	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.390	152	2042	0.365	ng/u1	0.00
18) Fluoranthene-d10	19.394	212	4634	0.363	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.397	88	798	0.413	ng/u1	99
5) Naphthalene	10.856	128	3629	0.365	ng/u1	99
7) 2-Methylnaphthalene	12.462	142	2229	0.364	ng/u1	99
8) 1-Methylnaphthalene	12.682	142	2271	0.365	ng/u1	98
10) Acenaphthylene	14.347	152	3448	0.350	ng/u1	99
11) Acenaphthene	14.689	153	2642	0.362	ng/u1	96
12) Fluorene	15.674	166	2958	0.357	ng/u1	99
14) Pentachlorophenol	17.029	266	605	0.382	ng/u1	96
15) Phenanthrene	17.413	178	4566	0.348	ng/u1	99
16) Anthracene	17.506	178	4290	0.355	ng/u1	98
19) Fluoranthene	19.422	202	5762	0.370	ng/u1	99
20) Pyrene	19.784	202	6017	0.362	ng/u1	99
21) Benzo(a)anthracene	21.527	228	5450	0.364	ng/u1	99
22) Chrysene	21.582	228	5892	0.373	ng/u1	99
24) Benzo(b)fluoranthene	23.228	252	6644	0.382	ng/u1	94
25) Benzo(k)fluoranthene	23.278	252	6635	0.383	ng/u1	94
26) Benzo(a)pyrene	23.865	252	6111	0.372	ng/u1	94
27) Indeno(1,2,3-cd)pyrene	26.501	276	8138	0.339	ng/u1#	60
28) Dibenzo(a,h)anthracene	26.525	278	6443	0.342	ng/u1	95
29) Benzo(g,h,i)perylene	27.290	276	7025	0.329	ng/u1	99

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

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