

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112122\
 Data File : BM037680.D
 Acq On : 23 Nov 2022 15:20
 Operator : CG/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4180

Quant Time: Nov 24 01:49:52 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM11622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 21 22:36:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.125	152	5276	0.400	ng/ul	0.00
4) Naphthalene-d8	10.942	136	15494	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.742	164	9051	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.475	188	20442	0.400	ng/ul	0.00
17) Chrysene-d12	21.636	240	14583	0.400	ng/ul	# 0.00
23) Perylene-d12	24.124	264	13137	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.450	96	2897	0.472	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.520	152	10020	0.390	ng/ul	0.00
18) Fluoranthene-d10	19.490	212	21403	0.444	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.484	88	2942	0.461	ng/ul#	58
5) Naphthalene	10.991	128	20003	0.425	ng/ul	98
7) 2-Methylnaphthalene	12.592	142	12868	0.402	ng/ul	100
8) 1-Methylnaphthalene	12.806	142	13104	0.400	ng/ul	99
10) Acenaphthylene	14.465	152	19255	0.383	ng/ul	100
11) Acenaphthene	14.803	153	15381	0.438	ng/ul	98
12) Fluorene	15.783	166	18015	0.426	ng/ul	99
14) Pentachlorophenol	17.120	266	2060	0.449	ng/ul	98
15) Phenanthrene	17.517	178	29857	0.427	ng/ul	100
16) Anthracene	17.610	178	26018	0.386	ng/ul	99
19) Fluoranthene	19.518	202	33171	0.471	ng/ul#	93
20) Pyrene	19.880	202	33941	0.464	ng/ul	97
21) Benzo(a)anthracene	21.619	228	27786	0.426	ng/ul	100
22) Chrysene	21.674	228	30151	0.466	ng/ul	100
24) Benzo(b)fluoranthene	23.358	252	31818	0.531	ng/ul	91
25) Benzo(k)fluoranthene	23.408	252	28620	0.473	ng/ul#	94
26) Benzo(a)pyrene	24.010	252	26574	0.510	ng/ul#	83
27) Indeno(1,2,3-cd)pyrene	26.721	276	33591	0.487	ng/ul#	86
28) Dibenzo(a,h)anthracene	26.741	278	25713	0.474	ng/ul#	95
29) Benzo(g,h,i)perylene	27.526	276	27390	0.468	ng/ul	95

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

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