

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030923\
 Data File : BM038940.D
 Acq On : 09 Mar 2023 11:37
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4075

Quant Time: Mar 09 13:36:19 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM030323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 13:36:03 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.995	152	12397	0.400	ng/u1	0.00
4) Naphthalene-d8	10.806	136	39364	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.633	164	21443	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.375	188	43383	0.400	ng/u1	0.00
17) Chrysene-d12	21.553	240	34848	0.400	ng/u1	0.00
23) Perylene-d12	23.999	264	25037	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.380	96	5884	0.419	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.389	152	19051	0.400	ng/u1	0.00
18) Fluoranthene-d10	19.398	212	33489	0.404	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.413	88	6135	0.432	ng/u1	97
5) Naphthalene	10.855	128	36767	0.391	ng/u1	100
7) 2-Methylnaphthalene	12.466	142	22577	0.395	ng/u1	100
8) 1-Methylnaphthalene	12.681	142	23476	0.392	ng/u1	99
10) Acenaphthylene	14.351	152	28976	0.372	ng/u1	99
11) Acenaphthene	14.693	153	24657	0.385	ng/u1	98
12) Fluorene	15.678	166	27570	0.381	ng/u1	100
14) Pentachlorophenol	17.020	266	3416	0.327	ng/u1	99
15) Phenanthrene	17.417	178	44103	0.385	ng/u1	99
16) Anthracene	17.505	178	36888	0.391	ng/u1	99
19) Fluoranthene	19.426	202	47079	0.399	ng/u1	98
20) Pyrene	19.789	202	48418	0.400	ng/u1	100
21) Benzo(a)anthracene	21.538	228	38185	0.387	ng/u1	99
22) Chrysene	21.591	228	41372	0.349	ng/u1	100
24) Benzo(b)fluoranthene	23.251	252	38435	0.393	ng/u1	98
25) Benzo(k)fluoranthene	23.301	252	36290	0.371	ng/u1	98
26) Benzo(a)pyrene	23.888	252	29735	0.381	ng/u1	97
27) Indeno(1,2,3-cd)pyrene	26.545	276	38095	0.386	ng/u1	98
28) Dibenzo(a,h)anthracene	26.568	278	30585	0.407	ng/u1	98
29) Benzo(g,h,i)perylene	27.326	276	33019	0.359	ng/u1	99

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

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