

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101921\
 Data File : BM032609.D
 Acq On : 20 Oct 2021 06:49
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4145

Quant Time: Oct 20 07:22:50 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 19 15:20:30 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.557	152	6395	0.40	ng/ul	0.00
4) Naphthalene-d8	10.344	136	19752	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.213	164	8872	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.946	188	15685	0.40	ng/ul	# 0.00
17) Chrysene-d12	21.123	240	11442	0.40	ng/ul	0.00
23) Perylene-d12	23.259	264	10593	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.927	96	3129	0.44	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.940	152	9787	0.37	ng/ul	0.00
18) Fluoranthene-d10	18.973	212	14316	0.40	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.962	88	3485	0.44	ng/ul#	82
5) Naphthalene	10.389	128	23644	0.39	ng/ul	99
7) 2-Methylnaphthalene	12.012	142	14588	0.37	ng/ul	98
8) 1-Methylnaphthalene	12.233	142	14254	0.37	ng/ul	99
10) Acenaphthylene	13.933	152	14295	0.38	ng/ul	98
11) Acenaphthene	14.274	153	14086	0.41	ng/ul	99
12) Fluorene	15.260	166	14874	0.38	ng/ul	98
14) Pentachlorophenol	16.620	266	1936	0.48	ng/ul	99
15) Phenanthrene	16.988	178	21464	0.41	ng/ul	100
16) Anthracene	17.078	178	17455	0.38	ng/ul	100
19) Fluoranthene	19.004	202	22472	0.40	ng/ul	99
20) Pyrene	19.366	202	22401	0.38	ng/ul	99
21) Benzo(a)anthracene	21.106	228	17074	0.38	ng/ul	100
22) Chrysene	21.157	228	20176	0.41	ng/ul	100
24) Benzo(b)fluoranthene	22.624	252	20551	0.42	ng/ul	100
25) Benzo(k)fluoranthene	22.665	252	20261	0.42	ng/ul	100
26) Benzo(a)pyrene	23.167	252	16838	0.41	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.369	276	22327	0.45	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.374	278	17719	0.45	ng/ul	99
29) Benzo(g,h,i)perylene	26.019	276	19920	0.46	ng/ul	99

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

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