

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031224\
 Data File : BM044584.D
 Acq On : 12 Mar 2024 09:29
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4052

Quant Time: Mar 12 10:35:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM031124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 11 14:02:17 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.724	152	6251	0.400	ng/ul	0.00
4) Naphthalene-d8	10.501	136	20532	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.363	164	11622	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.120	188	26048	0.400	ng/ul	0.00
17) Chrysene-d12	21.315	240	16501	0.400	ng/ul	0.00
23) Perylene-d12	23.574	264	15774	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.285	96	3209	0.392	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.101	152	11628	0.394	ng/ul	0.00
18) Fluoranthene-d10	19.151	212	23670	0.409	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.314	88	3041	0.363	ng/ul	90
5) Naphthalene	10.550	128	22329	0.387	ng/ul	100
7) 2-Methylnaphthalene	12.173	142	14327	0.392	ng/ul	99
8) 1-Methylnaphthalene	12.393	142	14493	0.393	ng/ul	99
10) Acenaphthylene	14.081	152	22932	0.388	ng/ul	99
11) Acenaphthene	14.428	153	16791	0.394	ng/ul	99
12) Fluorene	15.422	166	19344	0.397	ng/ul	98
14) Pentachlorophenol	16.774	266	2075	0.330	ng/ul	95
15) Phenanthrene	17.162	178	30491	0.390	ng/ul	99
16) Anthracene	17.255	178	28346	0.386	ng/ul	99
19) Fluoranthene	19.183	202	32795	0.406	ng/ul	99
20) Pyrene	19.546	202	33280	0.402	ng/ul	98
21) Benzo(a)anthracene	21.303	228	24404	0.388	ng/ul	99
22) Chrysene	21.353	228	27839	0.395	ng/ul	99
24) Benzo(b)fluoranthene	22.896	252	23991	0.398	ng/ul	100
25) Benzo(k)fluoranthene	22.940	252	25539	0.392	ng/ul	99
26) Benzo(a)pyrene	23.481	252	21012	0.384	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.862	276	23849	0.379	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.893	278	18032	0.372	ng/ul	100
29) Benzo(g,h,i)perylene	26.550	276	21567	0.391	ng/ul	100

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

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