

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072221\
 Data File : BM031051.D
 Acq On : 21 Jul 2021 20:23
 Operator : CG/JU
 Sample : SSTD0.409
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4010

Quant Time: Jul 22 01:41:51 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 01:40:36 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.642	152	931	0.400	ng/ul	0.00
4) Naphthalene-d8	10.406	136	3519	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.265	164	2303	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.013	188	5270	0.400	ng/ul	0.00
17) Chrysene-d12	21.200	240	5055	0.400	ng/ul	0.00
23) Perylene-d12	23.375	264	4737	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	386	0.403	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.998	152	2307	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.043	212	5777	0.340	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.251	88	339	0.317	ng/ul	100
5) Naphthalene	10.455	128	3966	0.378	ng/ul	100
7) 2-Methylnaphthalene	12.070	142	2871	0.387	ng/ul	100
8) 1-Methylnaphthalene	12.290	142	2764	0.381	ng/ul	100
10) Acenaphthylene	13.986	152	3589	0.318	ng/ul	100
11) Acenaphthene	14.330	153	3152	0.387	ng/ul	100
12) Fluorene	15.319	166	3705	0.383	ng/ul	100
14) Pentachlorophenol	16.662	266	642	0.345	ng/ul	100
15) Phenanthrene	17.055	178	6298	0.381	ng/ul	100
16) Anthracene	17.145	178	5881	0.371	ng/ul	100
19) Fluoranthene	19.073	202	7543	0.339	ng/ul	100
20) Pyrene	19.436	202	7723	0.343	ng/ul	100
21) Benzo(a)anthracene	21.183	228	7532	0.345	ng/ul	100
22) Chrysene	21.234	228	8277	0.367	ng/ul	100
24) Benzo(b)fluoranthene	22.725	252	8436	0.368	ng/ul	100
25) Benzo(k)fluoranthene	22.769	252	8374	0.363	ng/ul	100
26) Benzo(a)pyrene	23.280	252	7326	0.368	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.533	276	9271	0.343	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.553	278	7456	0.342	ng/ul	100
29) Benzo(g,h,i)perylene	26.192	276	8045	0.351	ng/ul	100

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

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