

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM103124\
 Data File : BM048384.D
 Acq On : 01 Nov 2024 07:43
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4067

Quant Time: Nov 03 08:06:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM103124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 31 14:58:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	3903	0.400	ng/ul	0.00
4) Naphthalene-d8	10.524	136	13391	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.363	164	7785	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.120	188	16363	0.400	ng/ul	0.00
17) Chrysene-d12	21.292	240	14213	0.400	ng/ul	0.00
23) Perylene-d12	23.531	264	14482	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.168	96	2244	0.468	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.141	152	6707	0.394	ng/ul	0.00
18) Fluoranthene-d10	19.137	212	15626	0.433	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.202	88	2534	0.487	ng/ul#	92
5) Naphthalene	10.574	128	13701	0.392	ng/ul	100
7) 2-Methylnaphthalene	12.212	142	8033	0.386	ng/ul	100
8) 1-Methylnaphthalene	12.410	142	8960	0.404	ng/ul	98
10) Acenaphthylene	14.086	152	12033	0.400	ng/ul	100
11) Acenaphthene	14.428	153	9745	0.404	ng/ul	99
12) Fluorene	15.432	166	10739	0.400	ng/ul	97
14) Pentachlorophenol	16.791	266	1339	0.294	ng/ul	91
15) Phenanthrene	17.158	178	15779	0.388	ng/ul	99
16) Anthracene	17.264	178	14426	0.379	ng/ul	99
19) Fluoranthene	19.169	202	20964	0.428	ng/ul	99
20) Pyrene	19.532	202	22443	0.427	ng/ul	100
21) Benzo(a)anthracene	21.280	228	16860	0.396	ng/ul	99
22) Chrysene	21.330	228	23999	0.410	ng/ul	100
24) Benzo(b)fluoranthene	22.855	252	18756	0.379	ng/ul	97
25) Benzo(k)fluoranthene	22.902	252	23284	0.397	ng/ul	98
26) Benzo(a)pyrene	23.440	252	18307	0.394	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.816	276	24777	0.381	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.839	278	19366	0.382	ng/ul	98
29) Benzo(g,h,i)perylene	26.500	276	21069	0.385	ng/ul	99

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

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