

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111324\
 Data File : BM048659.D
 Acq On : 14 Nov 2024 07:08
 Operator : RC/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4082

Quant Time: Nov 14 07:39:27 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM110624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 13 11:38:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.729	152	3709	0.400	ng/ul	0.00
4) Naphthalene-d8	10.484	136	13206	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.320	164	7693	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.081	188	15466	0.400	ng/ul	0.00
17) Chrysene-d12	21.260	240	12606	0.400	ng/ul	0.00
23) Perylene-d12	23.481	264	13662	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.126	96	2245	0.502	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.095	152	6496	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.108	212	14209	0.401	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.159	88	2361	0.459	ng/ul#	67
5) Naphthalene	10.534	128	13416	0.406	ng/ul	98
7) 2-Methylnaphthalene	12.172	142	8020	0.385	ng/ul	94
8) 1-Methylnaphthalene	12.370	142	9006	0.422	ng/ul	100
10) Acenaphthylene	14.043	152	12239	0.419	ng/ul	99
11) Acenaphthene	14.385	153	9755	0.417	ng/ul	99
12) Fluorene	15.389	166	10542	0.415	ng/ul	99
14) Pentachlorophenol	16.760	266	1333	0.235	ng/ul	98
15) Phenanthrene	17.123	178	15351	0.378	ng/ul	100
16) Anthracene	17.233	178	13845	0.366	ng/ul	99
19) Fluoranthene	19.135	202	19485	0.417	ng/ul	99
20) Pyrene	19.498	202	21021	0.415	ng/ul	99
21) Benzo(a)anthracene	21.251	228	14007	0.339	ng/ul	100
22) Chrysene	21.295	228	22022	0.445	ng/ul	99
24) Benzo(b)fluoranthene	22.812	252	18030	0.424	ng/ul	93
25) Benzo(k)fluoranthene	22.858	252	22064	0.442	ng/ul	93
26) Benzo(a)pyrene	23.396	252	17211	0.414	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	25.735	276	21918	0.378	ng/ul	98
28) Dibenzo(a,h)anthracene	25.752	278	16363	0.362	ng/ul	98
29) Benzo(g,h,i)perylene	26.416	276	19422	0.399	ng/ul	99

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

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