

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062424\
 Data File : BM046308.D
 Acq On : 26 Jun 2024 01:49
 Operator : MA/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 69 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4149

Quant Time: Jun 26 02:41:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM061824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 25 08:04:59 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.447	152	2034	0.400	ng/ul	0.00
4) Naphthalene-d8	10.222	136	6787	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.067	164	3811	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.837	188	7271	0.400	ng/ul	-0.02
17) Chrysene-d12	21.017	240	6024	0.400	ng/ul	0.00
23) Perylene-d12	23.104	264	8309	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.025	96	1047	0.375	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.844	152	3805	0.418	ng/ul	-0.03
18) Fluoranthene-d10	18.858	212	7332	0.406	ng/ul	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.054	88	1318	0.421	ng/ul#	79
5) Naphthalene	10.277	128	7567	0.400	ng/ul	98
7) 2-Methylnaphthalene	11.910	142	4393	0.412	ng/ul	99
8) 1-Methylnaphthalene	12.102	142	4709	0.403	ng/ul	95
10) Acenaphthylene	13.794	152	7997	0.442	ng/ul	98
11) Acenaphthene	14.132	153	5513	0.422	ng/ul	99
12) Fluorene	15.150	166	5879	0.420	ng/ul	97
14) Pentachlorophenol	16.516	266	894	0.410	ng/ul	98
15) Phenanthrene	16.875	178	8159	0.399	ng/ul	99
16) Anthracene	16.972	178	6537	0.430	ng/ul	98
19) Fluoranthene	18.886	202	11204	0.430	ng/ul	99
20) Pyrene	19.244	202	11680	0.414	ng/ul	99
21) Benzo(a)anthracene	21.003	228	9223	0.396	ng/ul	98
22) Chrysene	21.052	228	12107	0.408	ng/ul	98
24) Benzo(b)fluoranthene	22.487	252	10745	0.359	ng/ul	96
25) Benzo(k)fluoranthene	22.528	252	14646	0.412	ng/ul	99
26) Benzo(a)pyrene	23.013	252	11848	0.427	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.152	276	15537	0.342	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.158	278	12138	0.359	ng/ul	95
29) Benzo(g,h,i)perylene	25.769	276	12922	0.332	ng/ul	97

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

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