

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062223\
 Data File : BM040348.D
 Acq On : 22 Jun 2023 14:46
 Operator : MA/JU
 Sample : SST0.891
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.8025

Quant Time: Jun 23 00:38:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 23 00:36:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.953	152	6233	0.400	ng/ul	0.00
4) Naphthalene-d8	10.762	136	21403	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.591	164	10312	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.336	188	19105	0.400	ng/ul	0.00
17) Chrysene-d12	21.515	240	13855	0.400	ng/ul	0.00
23) Perylene-d12	23.932	264	14452	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.333	96	7674	0.784	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.351	152	23923	0.797	ng/ul	0.00
18) Fluoranthene-d10	19.361	212	35125	0.791	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.371	88	7860	0.789	ng/ul	93
5) Naphthalene	10.811	128	46158	0.782	ng/ul	99
7) 2-Methylnaphthalene	12.422	142	27802	0.802	ng/ul	100
8) 1-Methylnaphthalene	12.642	142	27790	0.802	ng/ul	100
10) Acenaphthylene	14.313	152	35912	0.780	ng/ul	99
11) Acenaphthene	14.656	153	30457	0.783	ng/ul	99
12) Fluorene	15.637	166	32909	0.792	ng/ul	99
14) Pentachlorophenol	16.999	266	2662	0.724	ng/ul	93
15) Phenanthrene	17.379	178	46513	0.784	ng/ul	98
16) Anthracene	17.467	178	40265	0.807	ng/ul	99
19) Fluoranthene	19.389	202	47067	0.800	ng/ul	99
20) Pyrene	19.751	202	48995	0.777	ng/ul	97
21) Benzo(a)anthracene	21.500	228	37289	0.796	ng/ul	99
22) Chrysene	21.553	228	42889	0.802	ng/ul	99
24) Benzo(b)fluoranthene	23.192	252	45006	0.805	ng/ul	93
25) Benzo(k)fluoranthene	23.239	252	44249	0.803	ng/ul	94
26) Benzo(a)pyrene	23.824	252	39744	0.789	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.437	276	56271	0.802	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.454	278	44527	0.811	ng/ul	96
29) Benzo(g,h,i)perylene	27.205	276	49001	0.797	ng/ul	97

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

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