

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071923\
 Data File : BM040964.D
 Acq On : 19 Jul 2023 20:14
 Operator : MA/JU
 Sample : SST0.496
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.4031

Quant Time: Jul 20 05:00:19 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 20 04:54:10 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.878	152	3361	0.400	ng/ul	0.00
4) Naphthalene-d8	10.680	136	13141	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.523	164	5078	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.282	188	9501	0.400	ng/ul	0.00
17) Chrysene-d12	21.472	240	6396	0.400	ng/ul	0.00
23) Perylene-d12	23.871	264	5737	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	1770	0.335	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.280	152	7417	0.403	ng/ul	0.00
18) Fluoranthene-d10	19.315	212	9381	0.458	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.317	88	1825	0.340	ng/ul	100
5) Naphthalene	10.730	128	14187	0.391	ng/ul	100
7) 2-Methylnaphthalene	12.352	142	8225	0.386	ng/ul	100
8) 1-Methylnaphthalene	12.566	142	8746	0.411	ng/ul	100
10) Acenaphthylene	14.250	152	8422	0.371	ng/ul	100
11) Acenaphthene	14.587	153	7433	0.388	ng/ul	100
12) Fluorene	15.582	166	7663	0.375	ng/ul	100
14) Pentachlorophenol	16.974	266	681	0.372	ng/ul	98
15) Phenanthrene	17.324	178	11183	0.379	ng/ul	100
16) Anthracene	17.421	178	8527	0.344	ng/ul	100
19) Fluoranthene	19.343	202	11903	0.438	ng/ul	100
20) Pyrene	19.705	202	12698	0.436	ng/ul	100
21) Benzo(a)anthracene	21.457	228	7409	0.342	ng/ul	100
22) Chrysene	21.507	228	8740	0.354	ng/ul	100
24) Benzo(b)fluoranthene	23.135	252	8380	0.377	ng/ul	100
25) Benzo(k)fluoranthene	23.187	252	8322	0.380	ng/ul	100
26) Benzo(a)pyrene	23.763	252	7309	0.365	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.351	276	9870	0.354	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.367	278	7476	0.343	ng/ul	100
29) Benzo(g,h,i)perylene	27.102	276	9301	0.381	ng/ul	100

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

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