

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013124\
 Data File : BM043932.D
 Acq On : 31 Jan 2024 12:56
 Operator : MA/JU
 Sample : SSTD0.498
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD0.4003

Quant Time: Jan 31 22:44:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM013124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 15:16:31 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.926	152	8648	0.400	ng/ul	0.00
4) Naphthalene-d8	10.727	136	23673	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.566	164	10983	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.318	188	23377	0.400	ng/ul	0.00
17) Chrysene-d12	21.517	240	17173	0.400	ng/ul	0.00
23) Perylene-d12	23.929	264	18821	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.365	96	4080	0.367	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.316	152	11270	0.363	ng/ul	0.00
18) Fluoranthene-d10	19.350	212	19359	0.359	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.403	88	4389	0.376	ng/ul	100
5) Naphthalene	10.776	128	24951	0.368	ng/ul	100
7) 2-Methylnaphthalene	12.393	142	14725	0.364	ng/ul	100
8) 1-Methylnaphthalene	12.607	142	14595	0.364	ng/ul	100
10) Acenaphthylene	14.288	152	21761	0.365	ng/ul	100
11) Acenaphthene	14.626	153	15032	0.364	ng/ul	100
12) Fluorene	15.616	166	16359	0.361	ng/ul	100
14) Pentachlorophenol	16.959	266	1800	0.325	ng/ul	100
15) Phenanthrene	17.360	178	25509	0.358	ng/ul	100
16) Anthracene	17.453	178	24121	0.357	ng/ul	100
19) Fluoranthene	19.382	202	28420	0.359	ng/ul	100
20) Pyrene	19.745	202	29129	0.358	ng/ul	100
21) Benzo(a)anthracene	21.500	228	25312	0.353	ng/ul	100
22) Chrysene	21.552	228	26858	0.358	ng/ul	100
24) Benzo(b)fluoranthene	23.192	252	26224	0.364	ng/ul	100
25) Benzo(k)fluoranthene	23.242	252	27276	0.358	ng/ul	100
26) Benzo(a)pyrene	23.821	252	23426	0.357	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.427	276	30015	0.356	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.457	278	23805	0.353	ng/ul	100
29) Benzo(g,h,i)perylene	27.191	276	26490	0.363	ng/ul	100

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

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