

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN031524\
 Data File : BN030406.D
 Acq On : 14 Mar 2024 21:01
 Operator : MA/JU
 Sample : SSTD0.871
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.8247

Quant Time: Mar 15 00:57:21 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN031524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 15 00:54:55 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.604	152	1683	0.400	ng/ul	0.00
4) Naphthalene-d8	10.377	136	3737	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.254	164	2170	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.009	188	4474	0.400	ng/ul	0.00
17) Chrysene-d12	21.215	240	3995	0.400	ng/ul	0.00
23) Perylene-d12	23.425	264	4587	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1655	0.710	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.978	152	4605	0.953	ng/ul	0.00
18) Fluoranthene-d10	19.046	212	9947	0.764	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.187	88	1734	0.686	ng/ul	95
5) Naphthalene	10.427	128	8231	0.746	ng/ul	95
7) 2-Methylnaphthalene	12.055	142	5200	0.870	ng/ul#	100
8) 1-Methylnaphthalene	12.275	142	5351	0.845	ng/ul#	28
10) Acenaphthylene	13.972	152	8434	0.695	ng/ul	98
11) Acenaphthene	14.319	153	6114	0.687	ng/ul	98
12) Fluorene	15.313	166	7010	0.773	ng/ul	98
14) Pentachlorophenol	16.662	266	1040	0.815	ng/ul#	100
15) Phenanthrene	17.051	178	10616	0.708	ng/ul	97
16) Anthracene	17.144	178	9750	0.739	ng/ul	97
19) Fluoranthene	19.079	202	12847	0.724	ng/ul	97
20) Pyrene	19.441	202	13519	0.701	ng/ul	96
21) Benzo(a)anthracene	21.198	228	11205	0.794	ng/ul	99
22) Chrysene	21.250	228	12634	0.745	ng/ul	97
24) Benzo(b)fluoranthene	22.761	252	12841	0.726	ng/ul	88
25) Benzo(k)fluoranthene	22.805	252	14806	0.710	ng/ul#	89
26) Benzo(a)pyrene	23.325	252	13039	0.724	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	25.631	276	17389	0.756	ng/ul#	90
28) Dibenzo(a,h)anthracene	25.658	278	13431	0.755	ng/ul#	87
29) Benzo(g,h,i)perylene	26.308	276	15324	0.718	ng/ul	96

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

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