

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011924\
 Data File : BM043879.D
 Acq On : 19 Jan 2024 19:20
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SICV049

Quant Time: Jan 19 22:49:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM011924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 19 22:45:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	3957	0.400	ng/ul	0.00
4) Naphthalene-d8	10.771	136	10384	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.584	164	4929	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.360	188	9330	0.400	ng/ul	0.00
17) Chrysene-d12	21.541	240	6126	0.400	ng/ul	0.00
23) Perylene-d12	23.961	264	7654	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	2215	0.411	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.376	152	5178	0.392	ng/ul	0.01
18) Fluoranthene-d10	19.378	212	7818	0.387	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.365	88	2259	0.388	ng/ul	98
5) Naphthalene	10.826	128	11964	0.396	ng/ul	99
7) 2-Methylnaphthalene	12.453	142	7078	0.404	ng/ul	100
8) 1-Methylnaphthalene	12.651	142	6793	0.377	ng/ul	99
10) Acenaphthylene	14.316	152	10719	0.407	ng/ul	99
11) Acenaphthene	14.649	153	7824	0.399	ng/ul	99
12) Fluorene	15.653	166	8274	0.408	ng/ul	100
14) Pentachlorophenol	17.149	266	353	0.298	ng/ul#	85
15) Phenanthrene	17.402	178	11505	0.389	ng/ul	99
16) Anthracene	17.508	178	10944	0.397	ng/ul	99
19) Fluoranthene	19.405	202	12799	0.393	ng/ul	97
20) Pyrene	19.773	202	13666	0.401	ng/ul	99
21) Benzo(a)anthracene	21.535	228	8634	0.363	ng/ul	100
22) Chrysene	21.582	228	13927	0.435	ng/ul	99
24) Benzo(b)fluoranthene	23.221	252	11568	0.397	ng/ul	97
25) Benzo(k)fluoranthene	23.274	252	13910	0.398	ng/ul	97
26) Benzo(a)pyrene	23.862	252	11429	0.384	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.484	276	15507	0.404	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.527	278	11428	0.390	ng/ul	96
29) Benzo(g,h,i)perylene	27.249	276	13559	0.396	ng/ul	98

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

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