

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100523\
 Data File : BN028065.D
 Acq On : 06 Oct 2023 01:13
 Operator : MA/JU
 Sample : PB156001BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS001

Quant Time: Oct 06 04:27:01 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 03 15:50:04 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	4953	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.757	136	15557	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.583	164	8039	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.337	188	15738	0.400	ng/ul	-0.01
17) Chrysene-d12	21.519	240	14474	0.400	ng/ul	#-0.02
23) Perylene-d12	23.980	264	16724	0.400	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	3715	0.597	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.341	152	6977	0.278	ng/ul	-0.01
18) Fluoranthene-d10	19.362	212	13936	0.304	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.363	88	9773	1.484	ng/ul	89
5) Naphthalene	10.806	128	13325	0.302	ng/ul	100
7) 2-Methylnaphthalene	12.412	142	7971	0.286	ng/ul	99
8) 1-Methylnaphthalene	12.632	142	8223	0.285	ng/ul	98
10) Acenaphthylene	14.310	152	10942	0.300	ng/ul	98
11) Acenaphthene	14.643	153	8637	0.298	ng/ul	99
12) Fluorene	15.633	166	9337	0.289	ng/ul	99
14) Pentachlorophenol	16.974	266	2245	0.665	ng/ul	98
15) Phenanthrene	17.380	178	14201	0.303	ng/ul	99
16) Anthracene	17.472	178	13409	0.320	ng/ul	98
19) Fluoranthene	19.394	202	17104	0.309	ng/ul	99
20) Pyrene	19.757	202	18263	0.319	ng/ul	98
21) Benzo(a)anthracene	21.501	228	17704	0.335	ng/ul	97
22) Chrysene	21.557	228	17447	0.327	ng/ul	97
24) Benzo(b)fluoranthene	23.217	252	18844	0.304	ng/ul#	62
25) Benzo(k)fluoranthene	23.267	252	19251	0.312	ng/ul#	68
26) Benzo(a)pyrene	23.872	252	16764	0.317	ng/ul#	63
27) Indeno(1,2,3-cd)pyrene	26.566	276	22958	0.359	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.586	278	18706	0.359	ng/ul#	75
29) Benzo(g,h,i)perylene	27.377	276	18655	0.357	ng/ul#	85

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

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