

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100523\
 Data File : BN028064.D
 Acq On : 06 Oct 2023 00:36
 Operator : MA/JU
 Sample : PB155952BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS952

Quant Time: Oct 06 04:26:51 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	7.941	152	5272	0.400	ng/ul	0.00
4) Naphthalene-d8	10.757	136	15985	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.583	164	8386	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.337	188	16878	0.400	ng/ul	-0.01
17) Chrysene-d12	21.519	240	15316	0.400	ng/ul	#-0.02
23) Perylene-d12	23.980	264	17621	0.400	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.330	96	3625	0.547	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.335	152	6945	0.270	ng/ul	-0.02
18) Fluoranthene-d10	19.362	212	14302	0.294	ng/ul	0.00
Target Compounds				Qvalue		
2) 1,4-Dioxane	3.364	88	9435	1.346	ng/ul	90
5) Naphthalene	10.806	128	12966	0.286	ng/ul	99
7) 2-Methylnaphthalene	12.412	142	7833	0.273	ng/ul	100
8) 1-Methylnaphthalene	12.632	142	8033	0.271	ng/ul	99
10) Acenaphthylene	14.310	152	11085	0.291	ng/ul	99
11) Acenaphthene	14.643	153	8526	0.282	ng/ul	99
12) Fluorene	15.633	166	9459	0.281	ng/ul	98
14) Pentachlorophenol	16.974	266	2403	0.664	ng/ul	98
15) Phenanthrene	17.380	178	14443	0.287	ng/ul	97
16) Anthracene	17.473	178	14032	0.312	ng/ul	98
19) Fluoranthene	19.394	202	17648	0.302	ng/ul	100
20) Pyrene	19.757	202	18468	0.305	ng/ul	97
21) Benzo(a)anthracene	21.501	228	17575	0.314	ng/ul	95
22) Chrysene	21.557	228	17332	0.307	ng/ul#	94
24) Benzo(b)fluoranthene	23.217	252	18806	0.288	ng/ul#	34
25) Benzo(k)fluoranthene	23.267	252	19076	0.293	ng/ul#	34
26) Benzo(a)pyrene	23.869	252	17142	0.308	ng/ul#	22
27) Indeno(1,2,3-cd)pyrene	26.563	276	23502	0.349	ng/ul#	90
28) Dibenzo(a,h)anthracene	26.583	278	19278	0.352	ng/ul#	53
29) Benzo(g,h,i)perylene	27.374	276	18640	0.338	ng/ul#	72

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

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