

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090523\
 Data File : BN027401.D
 Acq On : 06 Sep 2023 08:26
 Operator : MA/JU
 Sample : PB155063BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS063

Quant Time: Sep 06 08:57:54 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 02 06:16:48 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.055	152	4196	0.400	ng/ul	0.00
4) Naphthalene-d8	10.878	136	13557	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.685	164	7672	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.434	188	18470	0.400	ng/ul	0.00
17) Chrysene-d12	21.612	240	17763	0.400	ng/ul	0.00
23) Perylene-d12	24.135	264	16072	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.418	96	3365	0.603	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.456	152	6896	0.335	ng/ul	0.00
18) Fluoranthene-d10	19.455	212	19435	0.319	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.452	88	7807	1.382	ng/ul#	87
5) Naphthalene	10.928	128	11662	0.297	ng/ul	99
7) 2-Methylnaphthalene	12.528	142	7051	0.287	ng/ul	100
8) 1-Methylnaphthalene	12.742	142	7880	0.301	ng/ul	100
10) Acenaphthylene	14.416	152	10541	0.300	ng/ul	100
11) Acenaphthene	14.749	153	8924	0.295	ng/ul	99
12) Fluorene	15.735	166	10094	0.298	ng/ul	100
14) Pentachlorophenol	17.063	266	2469	0.686	ng/ul	100
15) Phenanthrene	17.476	178	17596	0.292	ng/ul	100
16) Anthracene	17.569	178	14221	0.297	ng/ul	100
19) Fluoranthene	19.483	202	22064	0.270	ng/ul	99
20) Pyrene	19.850	202	23005	0.277	ng/ul	98
21) Benzo(a)anthracene	21.597	228	19119	0.299	ng/ul	99
22) Chrysene	21.653	228	21532	0.288	ng/ul	99
24) Benzo(b)fluoranthene	23.351	252	21292	0.281	ng/ul	99
25) Benzo(k)fluoranthene	23.404	252	21474	0.289	ng/ul	98
26) Benzo(a)pyrene	24.024	252	16896	0.285	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.807	276	19343	0.294	ng/ul	99
28) Dibenzo(a,h)anthracene	26.834	278	14752	0.293	ng/ul	99
29) Benzo(g,h,i)perylene	27.629	276	17996	0.303	ng/ul	99

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

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