

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100523\
 Data File : BN028143.D
 Acq On : 08 Oct 2023 04:44
 Operator : MA/JU
 Sample : PB155976BS
 Misc :
 ALS Vial : 71 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS976

Quant Time: Oct 08 05:12:09 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 : Fri Oct 06 05:15:40 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	5179	0.400	ng/ul	0.00
4) Naphthalene-d8	10.746	136	15852	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.573	164	7736	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.329	188	14654	0.400	ng/ul	0.00
17) Chrysene-d12	21.516	240	12946	0.400	ng/ul	0.00
23) Perylene-d12	23.980	264	14044	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.321	96	4675	0.719	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.330	152	8050	0.315	ng/ul	0.00
18) Fluoranthene-d10	19.357	212	14479	0.353	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.355	88	12008	1.743	ng/ul#	84
5) Naphthalene	10.795	128	15331	0.341	ng/ul	100
7) 2-Methylnaphthalene	12.401	142	8945	0.315	ng/ul	99
8) 1-Methylnaphthalene	12.621	142	9476	0.322	ng/ul	99
10) Acenaphthylene	14.300	152	11577	0.330	ng/ul	99
11) Acenaphthene	14.638	153	9481	0.339	ng/ul	99
12) Fluorene	15.628	166	10172	0.327	ng/ul	98
14) Pentachlorophenol	16.974	266	1658	0.528	ng/ul	99
15) Phenanthrene	17.371	178	15404	0.353	ng/ul	99
16) Anthracene	17.468	178	13284	0.340	ng/ul	98
19) Fluoranthene	19.390	202	17221	0.348	ng/ul	100
20) Pyrene	19.752	202	18026	0.352	ng/ul	98
21) Benzo(a)anthracene	21.501	228	16749	0.354	ng/ul	100
22) Chrysene	21.554	228	17526	0.367	ng/ul	99
24) Benzo(b)fluoranthene	23.214	252	18737	0.360	ng/ul	98
25) Benzo(k)fluoranthene	23.263	252	19069	0.368	ng/ul	98
26) Benzo(a)pyrene	23.869	252	15945	0.359	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.569	276	22311	0.415	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.586	278	18264	0.418	ng/ul	99
29) Benzo(g,h,i)perylene	27.374	276	18857	0.429	ng/ul	98

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

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