

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091223\
 Data File : BN027583.D
 Acq On : 11 Sep 2023 19:43
 Operator : MA/JU
 Sample : PB155158BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS158

Quant Time: Sep 12 02:13:49 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 : Tue Sep 12 02:09:21 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.042	152	6803	0.400	ng/ul	0.00
4) Naphthalene-d8	10.861	136	18243	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.666	164	8123	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.418	188	15763	0.400	ng/ul	0.00
17) Chrysene-d12	21.592	240	12193	0.400	ng/ul	0.00
23) Perylene-d12	24.097	264	11488	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.439	96	4220	0.466	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.434	152	7038	0.254	ng/ul	0.00
18) Fluoranthene-d10	19.436	212	11278	0.270	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.473	88	11876	1.297	ng/ul#	88
5) Naphthalene	10.911	128	14827	0.280	ng/ul	100
7) 2-Methylnaphthalene	12.506	142	8472	0.256	ng/ul	99
8) 1-Methylnaphthalene	12.726	142	8614	0.244	ng/ul	100
10) Acenaphthylene	14.398	152	10029	0.270	ng/ul	100
11) Acenaphthene	14.731	153	8541	0.267	ng/ul	98
12) Fluorene	15.716	166	9061	0.253	ng/ul	99
14) Pentachlorophenol	17.050	266	1786	0.581	ng/ul	99
15) Phenanthrene	17.460	178	13913	0.271	ng/ul	100
16) Anthracene	17.553	178	11409	0.279	ng/ul	100
19) Fluoranthene	19.469	202	14953	0.267	ng/ul	99
20) Pyrene	19.831	202	15338	0.269	ng/ul	99
21) Benzo(a)anthracene	21.574	228	14113	0.321	ng/ul	99
22) Chrysene	21.633	228	14444	0.282	ng/ul	99
24) Benzo(b)fluoranthene	23.316	252	14464	0.267	ng/ul	94
25) Benzo(k)fluoranthene	23.369	252	14119	0.266	ng/ul	94
26) Benzo(a)pyrene	23.986	252	11975	0.283	ng/ul	93
27) Indeno(1,2,3-cd)pyrene	26.741	276	15906	0.338	ng/ul#	45
28) Dibenzo(a,h)anthracene	26.757	278	12812	0.356	ng/ul	99
29) Benzo(g,h,i)perylene	27.562	276	14077	0.331	ng/ul	98

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

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