

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071123\
 Data File : BN026462.D
 Acq On : 11 Jul 2023 14:40
 Operator : MA/JU
 Sample : PB153990BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS990

Quant Time: Jul 12 02:12:10 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 09 15:51:14 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	4842	0.400	ng/ul	0.00
4) Naphthalene-d8	10.732	136	17049	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.566	164	9036	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.322	188	13430	0.400	ng/ul	0.00
17) Chrysene-d12	21.523	240	8774	0.400	ng/ul	0.00
23) Perylene-d12	23.976	264	8823	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	3704	0.679	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.327	152	7524	0.287	ng/ul	0.00
18) Fluoranthene-d10	19.349	212	11482	0.344	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	8491	1.467	ng/ul#	87
5) Naphthalene	10.781	128	15804	0.323	ng/ul	99
7) 2-Methylnaphthalene	12.403	142	7813	0.251	ng/ul	98
8) 1-Methylnaphthalene	12.612	142	9175	0.285	ng/ul	100
10) Acenaphthylene	14.288	152	12098	0.286	ng/ul	99
11) Acenaphthene	14.626	153	10255	0.296	ng/ul	99
12) Fluorene	15.625	166	10028	0.253	ng/ul	98
14) Pentachlorophenol	16.988	266	2140	0.445	ng/ul	99
15) Phenanthrene	17.364	178	13272	0.299	ng/ul	98
16) Anthracene	17.465	178	10544	0.267	ng/ul	96
19) Fluoranthene	19.382	202	14297	0.305	ng/ul	98
20) Pyrene	19.744	202	15409	0.316	ng/ul	98
21) Benzo(a)anthracene	21.509	228	10150	0.280	ng/ul	99
22) Chrysene	21.558	228	11815	0.314	ng/ul	99
24) Benzo(b)fluoranthene	23.227	252	11612	0.308	ng/ul	89
25) Benzo(k)fluoranthene	23.277	252	12253	0.324	ng/ul	92
26) Benzo(a)pyrene	23.876	252	10033	0.307	ng/ul#	87
27) Indeno(1,2,3-cd)pyrene	26.558	276	11634	0.326	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.571	278	8758	0.324	ng/ul#	88
29) Benzo(g,h,i)perylene	27.339	276	10705	0.337	ng/ul	96

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

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