

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081222\
 Data File : BN021213.D
 Acq On : 13 Aug 2022 15:44
 Operator : CG/JU
 Sample : PB146961BS
 Misc :
 ALS Vial : 99 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS961

Quant Time: Aug 14 23:05:03 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 13 04:02:44 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.797	152	2154	0.400	ng/ul	0.00
4) Naphthalene-d8	10.590	136	7664	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.437	164	4472	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.196	188	7822	0.400	ng/ul	0.00
17) Chrysene-d12	21.398	240	5121	0.400	ng/ul	0.00
23) Perylene-d12	23.728	264	3979	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.194	96	2282	0.828	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.190	152	4358	0.388	ng/ul	0.00
18) Fluoranthene-d10	19.229	212	8755	0.554	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.228	88	2911	1.025	ng/ul#	93
5) Naphthalene	10.639	128	4038	0.172	ng/ul	97
7) 2-Methylnaphthalene	12.267	142	2315	0.162	ng/ul	100
8) 1-Methylnaphthalene	12.481	142	2657	0.186	ng/ul	99
10) Acenaphthylene	14.159	152	3510	0.168	ng/ul	97
11) Acenaphthene	14.502	153	2928	0.169	ng/ul	99
12) Fluorene	15.501	166	3068	0.162	ng/ul	99
14) Pentachlorophenol	16.875	266	206	0.171	ng/ul	97
15) Phenanthrene	17.234	178	4751	0.170	ng/ul	98
16) Anthracene	17.335	178	3601	0.166	ng/ul	97
19) Fluoranthene	19.262	202	5208	0.230	ng/ul	98
20) Pyrene	19.624	202	5413	0.231	ng/ul	95
21) Benzo(a)anthracene	21.384	228	2794	0.167	ng/ul	97
22) Chrysene	21.436	228	3938	0.179	ng/ul	98
24) Benzo(b)fluoranthene	23.023	252	3367	0.183	ng/ul	84
25) Benzo(k)fluoranthene	23.070	252	3563	0.183	ng/ul#	87
26) Benzo(a)pyrene	23.628	252	3180	0.189	ng/ul#	68
27) Indeno(1,2,3-cd)pyrene	26.146	276	3403	0.170	ng/ul	89
28) Dibenzo(a,h)anthracene	26.166	278	2590	0.163	ng/ul#	84
29) Benzo(g,h,i)perylene	26.874	276	3322	0.186	ng/ul	93

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

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