

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121324\
 Data File : BN035610.D
 Acq On : 13 Dec 2024 23:05
 Operator : RC/JU
 Sample : PB165482BS
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS482

Quant Time: Dec 13 23:28:08 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 10 16:10:57 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.292	152	2442	0.400	ng/ul	0.00
4) Naphthalene-d8	10.036	136	7417	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.953	164	4827	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.721	188	8882	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.964	240	6365	0.400	ng/ul	# 0.00
23) Perylene-d12	23.056	264	7313	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	1230	0.551	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.642	152	3634	0.334	ng/ul	0.00
18) Fluoranthene-d10	18.772	212	7137	0.406	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	2.993	88	3252	1.342	ng/ul#	87
5) Naphthalene	10.086	128	5535	0.296	ng/ul	98
7) 2-Methylnaphthalene	11.714	142	3844	0.297	ng/ul	97
8) 1-Methylnaphthalene	11.945	142	3921	0.300	ng/ul	98
10) Acenaphthylene	13.666	152	5830	0.304	ng/ul	98
11) Acenaphthene	14.013	153	4149	0.283	ng/ul	98
12) Fluorene	15.017	166	4737	0.275	ng/ul	98
14) Pentachlorophenol	16.388	266	1222	0.591	ng/ul	97
15) Phenanthrene	16.764	178	7097	0.302	ng/ul	98
16) Anthracene	16.852	178	6741	0.309	ng/ul	97
19) Fluoranthene	18.805	202	8037	0.363	ng/ul#	95
20) Pyrene	19.172	202	8505	0.369	ng/ul#	94
21) Benzo(a)anthracene	20.949	228	6951	0.315	ng/ul#	93
22) Chrysene	21.002	228	7192	0.324	ng/ul#	95
24) Benzo(b)fluoranthene	22.442	252	7187	0.297	ng/ul#	70
25) Benzo(k)fluoranthene	22.483	252	7721	0.310	ng/ul#	64
26) Benzo(a)pyrene	22.966	252	7009	0.317	ng/ul#	59
27) Indeno(1,2,3-cd)pyrene	25.087	276	9108	0.305	ng/ul#	1
28) Dibenzo(a,h)anthracene	25.104	278	7159	0.306	ng/ul#	69
29) Benzo(g,h,i)perylene	25.704	276	7433	0.316	ng/ul#	71

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

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