

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034934.D
 Acq On : 09 Nov 2024 10:27
 Operator : RC/JU
 Sample : PB164773BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS773

Quant Time: Nov 09 22:51:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.571	152	2502	0.400	ng/ul	0.00
4) Naphthalene-d8	10.333	136	6304	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.203	164	4849	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	12460	0.400	ng/ul	0.00
17) Chrysene-d12	21.145	240	12359	0.400	ng/ul	0.00
23) Perylene-d12	23.308	264	11514	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1604	0.674	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.934	152	3916	0.406	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	13491	0.425	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.182	88	2232	0.855	ng/ul	93
5) Naphthalene	10.383	128	2901	0.184	ng/ul	99
7) 2-Methylnaphthalene	12.005	142	2148	0.182	ng/ul	98
8) 1-Methylnaphthalene	12.225	142	2175	0.184	ng/ul	99
10) Acenaphthylene	13.921	152	3548	0.179	ng/ul	98
11) Acenaphthene	14.268	153	2731	0.182	ng/ul	98
12) Fluorene	15.258	166	3376	0.177	ng/ul	99
14) Pentachlorophenol	16.603	266	1022	0.355	ng/ul	96
15) Phenanthrene	16.992	178	6028	0.181	ng/ul	99
16) Anthracene	17.084	178	5413	0.173	ng/ul	99
19) Fluoranthene	19.014	202	7803	0.189	ng/ul#	97
20) Pyrene	19.376	202	8095	0.189	ng/ul#	97
21) Benzo(a)anthracene	21.128	228	8150	0.182	ng/ul	99
22) Chrysene	21.180	228	8596	0.188	ng/ul	99
24) Benzo(b)fluoranthene	22.665	252	8137	0.201	ng/ul	92
25) Benzo(k)fluoranthene	22.709	252	8027	0.189	ng/ul#	95
26) Benzo(a)pyrene	23.214	252	6836	0.189	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.463	276	7752	0.175	ng/ul#	96
28) Dibenzo(a,h)anthracene	25.480	278	6134	0.178	ng/ul#	95
29) Benzo(g,h,i)perylene	26.117	276	6191	0.183	ng/ul	97

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

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