

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034927.D
 Acq On : 09 Nov 2024 05:41
 Operator : RC/JU
 Sample : PB164771BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS771

Quant Time: Nov 09 06:03:17 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	2351	0.400	ng/ul	0.00
4) Naphthalene-d8	10.333	136	5960	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.208	164	4523	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.954	188	11560	0.400	ng/ul	0.00
17) Chrysene-d12	21.145	240	11222	0.400	ng/ul	0.00
23) Perylene-d12	23.314	264	10320	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1604	0.717	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.934	152	3904	0.428	ng/ul	0.00
18) Fluoranthene-d10	18.986	212	13083	0.454	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.187	88	2337	0.953	ng/ul#	94
5) Naphthalene	10.383	128	3101	0.208	ng/ul	98
7) 2-Methylnaphthalene	12.005	142	2299	0.207	ng/ul	99
8) 1-Methylnaphthalene	12.231	142	2297	0.205	ng/ul	99
10) Acenaphthylene	13.921	152	3688	0.200	ng/ul	99
11) Acenaphthene	14.268	153	2827	0.202	ng/ul	100
12) Fluorene	15.258	166	3533	0.198	ng/ul	99
14) Pentachlorophenol	16.608	266	1071	0.401	ng/ul	98
15) Phenanthrene	16.996	178	6233	0.202	ng/ul	99
16) Anthracene	17.085	178	5692	0.196	ng/ul	99
19) Fluoranthene	19.014	202	8135	0.217	ng/ul#	96
20) Pyrene	19.376	202	8340	0.214	ng/ul#	96
21) Benzo(a)anthracene	21.131	228	8279	0.204	ng/ul	100
22) Chrysene	21.180	228	8794	0.212	ng/ul	99
24) Benzo(b)fluoranthene	22.668	252	7978	0.220	ng/ul	95
25) Benzo(k)fluoranthene	22.712	252	8127	0.214	ng/ul#	95
26) Benzo(a)pyrene	23.217	252	6744	0.208	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	25.467	276	7716	0.195	ng/ul#	67
28) Dibenzo(a,h)anthracene	25.483	278	6028	0.195	ng/ul	94
29) Benzo(g,h,i)perylene	26.121	276	6145	0.203	ng/ul#	95

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

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