

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034991.D
 Acq On : 10 Nov 2024 22:17
 Operator : RC/JU
 Sample : PB164814BS
 Misc :
 ALS Vial : 85 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS814

Quant Time: Nov 10 23:32:44 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	7.562	152	2524	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.328	136	5981	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.199	164	4344	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	10387	0.400	ng/ul	0.00
17) Chrysene-d12	21.140	240	10324	0.400	ng/ul	0.00
23) Perylene-d12	23.306	264	11104	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	1967	0.819	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	4346	0.474	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	12790	0.482	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.183	88	5271	2.003	ng/ul#	92
5) Naphthalene	10.378	128	6420	0.429	ng/ul	99
7) 2-Methylnaphthalene	12.000	142	4608	0.412	ng/ul	99
8) 1-Methylnaphthalene	12.220	142	5387	0.480	ng/ul	99
10) Acenaphthylene	13.917	152	7225	0.407	ng/ul	100
11) Acenaphthene	14.264	153	5537	0.412	ng/ul	99
12) Fluorene	15.254	166	6755	0.395	ng/ul	99
14) Pentachlorophenol	16.599	266	2016	0.839	ng/ul	97
15) Phenanthrene	16.988	178	11642	0.419	ng/ul	99
16) Anthracene	17.081	178	10673	0.409	ng/ul	99
19) Fluoranthene	19.010	202	14724	0.427	ng/ul#	97
20) Pyrene	19.372	202	15306	0.428	ng/ul#	97
21) Benzo(a)anthracene	21.125	228	15494	0.414	ng/ul	99
22) Chrysene	21.175	228	16043	0.421	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	16263	0.417	ng/ul	100
25) Benzo(k)fluoranthene	22.703	252	17004	0.416	ng/ul	99
26) Benzo(a)pyrene	23.212	252	14529	0.417	ng/ul#	99
27) Indeno(1,2,3-cd)pyrene	25.454	276	19984	0.468	ng/ul	94
28) Dibenzo(a,h)anthracene	25.474	278	15623	0.470	ng/ul	98
29) Benzo(g,h,i)perylene	26.108	276	15487	0.475	ng/ul	98

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

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