

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020325\
 Data File : BN036253.D
 Acq On : 04 Feb 2025 06:14
 Operator : RC/JU
 Sample : PB166394BS
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS394

Quant Time: Feb 04 08:35:46 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN012125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 03 12:27:24 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.783	152	2942	0.400	ng/ul	0.00
4) Naphthalene-d8	10.578	136	7258	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.427	164	4208	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.174	188	8803	0.400	ng/ul	0.00
17) Chrysene-d12	21.366	240	8616	0.400	ng/ul	0.00
23) Perylene-d12	23.657	264	7530	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.256	96	1860	0.577	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.172	152	3720	0.406	ng/ul	0.00
18) Fluoranthene-d10	19.206	212	9319	0.394	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.290	88	5281	1.536	ng/ul	89
5) Naphthalene	10.627	128	7085	0.354	ng/ul	99
7) 2-Methylnaphthalene	12.244	142	4457	0.359	ng/ul	100
8) 1-Methylnaphthalene	12.464	142	4528	0.359	ng/ul	99
10) Acenaphthylene	14.144	152	6289	0.317	ng/ul	98
11) Acenaphthene	14.491	153	4997	0.342	ng/ul	98
12) Fluorene	15.481	166	5369	0.333	ng/ul	99
14) Pentachlorophenol	16.832	266	1628	0.528	ng/ul	99
15) Phenanthrene	17.216	178	8847	0.343	ng/ul	99
16) Anthracene	17.309	178	7801	0.326	ng/ul	99
19) Fluoranthene	19.234	202	11326	0.346	ng/ul	96
20) Pyrene	19.596	202	11874	0.347	ng/ul#	94
21) Benzo(a)anthracene	21.348	228	10632	0.346	ng/ul	100
22) Chrysene	21.398	228	11177	0.356	ng/ul	99
24) Benzo(b)fluoranthene	22.964	252	9715	0.387	ng/ul	89
25) Benzo(k)fluoranthene	23.011	252	9786	0.370	ng/ul#	91
26) Benzo(a)pyrene	23.555	252	8282	0.365	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	25.991	276	9519	0.317	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.004	278	7497	0.324	ng/ul	94
29) Benzo(g,h,i)perylene	26.705	276	7419	0.318	ng/ul	96

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

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