

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030112.D
 Acq On : 28 Feb 2024 03:22
 Operator : MA/JU
 Sample : PB159252BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS252

Quant Time: Feb 28 03:49:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	4279	0.400	ng/ul	0.00
4) Naphthalene-d8	10.649	136	11872	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	4727	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.262	188	9136	0.400	ng/ul	0.00
17) Chrysene-d12	21.461	240	6883	0.400	ng/ul	0.00
23) Perylene-d12	23.823	264	5933	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	3914	0.734	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.244	152	5075	0.336	ng/ul	0.00
18) Fluoranthene-d10	19.293	212	8378	0.370	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.349	88	10898	1.876	ng/ul#	81
5) Naphthalene	10.698	128	11642	0.348	ng/ul	99
7) 2-Methylnaphthalene	12.321	142	6210	0.324	ng/ul	99
8) 1-Methylnaphthalene	12.535	142	6586	0.338	ng/ul	99
10) Acenaphthylene	14.218	152	7831	0.340	ng/ul	98
11) Acenaphthene	14.561	153	6259	0.357	ng/ul	97
12) Fluorene	15.565	166	6691	0.365	ng/ul	98
14) Pentachlorophenol	16.916	266	1388	0.681	ng/ul#	49
15) Phenanthrene	17.304	178	10017	0.351	ng/ul	96
16) Anthracene	17.402	178	8325	0.347	ng/ul	96
19) Fluoranthene	19.326	202	11224	0.355	ng/ul	99
20) Pyrene	19.683	202	12235	0.378	ng/ul	99
21) Benzo(a)anthracene	21.447	228	7883	0.318	ng/ul	97
22) Chrysene	21.496	228	10021	0.369	ng/ul	99
24) Benzo(b)fluoranthene	23.104	252	8868	0.397	ng/ul	96
25) Benzo(k)fluoranthene	23.153	252	10635	0.430	ng/ul	97
26) Benzo(a)pyrene	23.721	252	8466	0.403	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.289	276	9586	0.376	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.329	278	7270	0.364	ng/ul#	89
29) Benzo(g,h,i)perylene	27.013	276	9105	0.394	ng/ul	98

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

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