

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020824\
 Data File : BN029595.D
 Acq On : 09 Feb 2024 02:11
 Operator : MA/JU
 Sample : P1303-04
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0FG9

Quant Time: Feb 09 02:53:24 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 : Wed Feb 07 02:17:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.884	152	5281	0.400	ng/ul	0.00
4) Naphthalene-d8	10.683	136	14257	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.530	164	6643	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.276	188	12717	0.400	ng/ul	0.00
17) Chrysene-d12	21.472	240	9389	0.400	ng/ul	0.00
23) Perylene-d12	23.845	264	11139	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	33760	5.127	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.278	152	5657	0.312	ng/ul	0.00
18) Fluoranthene-d10	19.309	212	9299	0.301	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.733	128	4305	0.107	ng/ul#	95
7) 2-Methylnaphthalene	12.350	142	1152	0.050	ng/ul	100
15) Phenanthrene	17.318	178	2934	0.074	ng/ul#	94
19) Fluoranthene	19.341	202	4108	0.095	ng/ul#	70
20) Pyrene	19.704	202	3480	0.079	ng/ul#	89
21) Benzo(a)anthracene	21.457	228	1413	0.042	ng/ul#	84
22) Chrysene	21.510	228	2204	0.059	ng/ul#	91
24) Benzo(b)fluoranthene	23.126	252	2871	0.069	ng/ul#	33
25) Benzo(k)fluoranthene	23.173	252	1455	0.031	ng/ul#	1
26) Benzo(a)pyrene	23.737	252	1417	0.036	ng/ul#	14
27) Indeno(1,2,3-cd)pyrene	26.297	276	1300	0.027	ng/ul#	100
29) Benzo(g,h,i)perylene	27.045	276	1398	0.032	ng/ul#	60

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

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