

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121322\
 Data File : BN023197.D
 Acq On : 14 Dec 2022 11:59
 Operator : CG/JU
 Sample : N5948-09MS
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EXZ94MS

Quant Time: Dec 14 23:19:41 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN112622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 13 23:52:47 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.015	152	5063	0.400 ng/u1	0.00
4) Naphthalene-d8	10.832	136	16530	0.400 ng/u1	0.00
9) Acenaphthene-d10	14.659	164	9882	0.400 ng/u1	0.00
13) Phenanthrene-d10	17.403	188	22656	0.400 ng/u1	0.00
17) Chrysene-d12	21.584	240	22096	0.400 ng/u1	0.00
23) Perylene-d12	24.039	264	18025	0.400 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.357	96	1087	0.179 ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.416	152	8379	0.333 ng/u1	0.00
18) Fluoranthene-d10	19.430	212	23366	0.286 ng/u1	0.00
Target Compounds					
					Qvalue
2) 1,4-Dioxane	3.395	88	3755	0.593 ng/u1	94
5) Naphthalene	10.881	128	68773	1.371 ng/u1	99
7) 2-Methylnaphthalene	12.493	142	13472	0.430 ng/u1	100
8) 1-Methylnaphthalene	12.707	142	14828	0.465 ng/u1	100
10) Acenaphthylene	14.377	152	13978	0.287 ng/u1	98
11) Acenaphthene	14.719	153	12120	0.314 ng/u1	97
12) Fluorene	15.705	166	14505	0.332 ng/u1	100
14) Pentachlorophenol	17.052	266	4117	0.869 ng/u1#	100
15) Phenanthrene	17.445	178	25967	0.334 ng/u1	99
16) Anthracene	17.534	178	21537	0.323 ng/u1	100
19) Fluoranthene	19.457	202	32570	0.285 ng/u1	97
20) Pyrene	19.820	202	33607	0.300 ng/u1	97
21) Benzo(a)anthracene	21.569	228	32429	0.384 ng/u1	99
22) Chrysene	21.622	228	33546	0.362 ng/u1	99
24) Benzo(b)fluoranthene	23.285	252	33929	0.371 ng/u1	96
25) Benzo(k)fluoranthene	23.335	252	31991	0.331 ng/u1	97
26) Benzo(a)pyrene	23.928	252	27203	0.391 ng/u1#	94
27) Indeno(1,2,3-cd)pyrene	26.604	276	34176	0.413 ng/u1#	100
28) Dibenzo(a,h)anthracene	26.621	278	26886	0.409 ng/u1	99
29) Benzo(g,h,i)perylene	27.392	276	27266	0.378 ng/u1	100

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

