

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030111.D
 Acq On : 28 Feb 2024 02:46
 Operator : MA/JU
 Sample : PB159257BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS257

Quant Time: Feb 28 03:19:43 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	4195	0.400	ng/ul	0.00
4) Naphthalene-d8	10.645	136	12175	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.497	164	5388	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.259	188	9412	0.400	ng/ul	0.00
17) Chrysene-d12	21.458	240	6732	0.400	ng/ul	0.00
23) Perylene-d12	23.817	264	5454	0.400	ng/ul #-	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	4033	0.771	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.245	152	5678	0.367	ng/ul	0.00
18) Fluoranthene-d10	19.290	212	9293	0.420	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.349	88	11303	1.985	ng/ul#	84
5) Naphthalene	10.694	128	12564	0.367	ng/ul	99
7) 2-Methylnaphthalene	12.317	142	6959	0.354	ng/ul	99
8) 1-Methylnaphthalene	12.537	142	7349	0.367	ng/ul	99
10) Acenaphthylene	14.220	152	9487	0.361	ng/ul	98
11) Acenaphthene	14.562	153	7716	0.386	ng/ul	99
12) Fluorene	15.561	166	7779	0.372	ng/ul	98
14) Pentachlorophenol	16.909	266	1674	0.797	ng/ul#	50
15) Phenanthrene	17.301	178	11117	0.378	ng/ul	97
16) Anthracene	17.399	178	9334	0.377	ng/ul	97
19) Fluoranthene	19.323	202	12547	0.406	ng/ul	100
20) Pyrene	19.680	202	13440	0.424	ng/ul	99
21) Benzo(a)anthracene	21.444	228	8444	0.348	ng/ul	99
22) Chrysene	21.493	228	10882	0.410	ng/ul	99
24) Benzo(b)fluoranthene	23.098	252	9185	0.448	ng/ul	96
25) Benzo(k)fluoranthene	23.147	252	11047	0.485	ng/ul	97
26) Benzo(a)pyrene	23.715	252	8681	0.450	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.282	276	9703	0.414	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.329	278	7343	0.400	ng/ul#	89
29) Benzo(g,h,i)perylene	27.006	276	9281	0.437	ng/ul	99

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

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