

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030185.D
 Acq On : 02 Mar 2024 17:31
 Operator : MA/JU
 Sample : PB159333BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS333

Quant Time: Mar 03 22:48:11 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN030224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 02 16:09:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.838	152	1375	0.400	ng/ul	0.00
4) Naphthalene-d8	10.634	136	3557	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.488	164	1632	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.247	188	2856	0.400	ng/ul	0.00
17) Chrysene-d12	21.456	240	1971	0.400	ng/ul	0.00
23) Perylene-d12	23.812	264	2007	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	1511	0.794	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.240	152	1898	0.413	ng/ul	0.00
18) Fluoranthene-d10	19.281	212	3164	0.493	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.340	88	4243	2.054	ng/ul#	85
5) Naphthalene	10.689	128	4336	0.413	ng/ul	99
7) 2-Methylnaphthalene	12.311	142	2270	0.399	ng/ul	99
8) 1-Methylnaphthalene	12.525	142	2455	0.407	ng/ul	98
10) Acenaphthylene	14.206	152	3699	0.405	ng/ul	97
11) Acenaphthene	14.548	153	2697	0.403	ng/ul	98
12) Fluorene	15.547	166	2793	0.410	ng/ul	97
14) Pentachlorophenol	16.897	266	753	0.924	ng/ul	98
15) Phenanthrene	17.289	178	3961	0.414	ng/ul	100
16) Anthracene	17.386	178	3427	0.407	ng/ul	98
19) Fluoranthene	19.314	202	4286	0.490	ng/ul	100
20) Pyrene	19.676	202	4610	0.484	ng/ul	98
21) Benzo(a)anthracene	21.438	228	3024	0.434	ng/ul	99
22) Chrysene	21.488	228	3732	0.446	ng/ul	98
24) Benzo(b)fluoranthene	23.098	252	3562	0.460	ng/ul	96
25) Benzo(k)fluoranthene	23.145	252	3967	0.435	ng/ul	97
26) Benzo(a)pyrene	23.712	252	3444	0.437	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.249	276	4376	0.435	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.282	278	3396	0.436	ng/ul	96
29) Benzo(g,h,i)perylene	26.983	276	4087	0.438	ng/ul	95

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

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