

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035305.D
 Acq On : 26 Nov 2024 06:56
 Operator : RC/JU
 Sample : PB165232BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS232

Quant Time: Nov 26 07:18:33 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 26 06:09:15 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.314	152	2937	0.400	ng/ul	0.00
4) Naphthalene-d8	10.064	136	7102	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.972	164	3820	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.738	188	7710	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.982	240	6344	0.400	ng/ul	# 0.00
23) Perylene-d12	23.080	264	6933	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1941	0.723	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.664	152	3420	0.329	ng/ul	0.00
18) Fluoranthene-d10	18.791	212	6845	0.391	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.005	88	5078	1.743	ng/ul	88
5) Naphthalene	10.108	128	6231	0.348	ng/ul	100
7) 2-Methylnaphthalene	11.741	142	3947	0.318	ng/ul	98
8) 1-Methylnaphthalene	11.967	142	4027	0.322	ng/ul	99
10) Acenaphthylene	13.685	152	5585	0.369	ng/ul	100
11) Acenaphthene	14.037	153	4259	0.367	ng/ul	98
12) Fluorene	15.036	166	4635	0.341	ng/ul	99
14) Pentachlorophenol	16.396	266	1211	0.674	ng/ul	99
15) Phenanthrene	16.781	178	7411	0.364	ng/ul	99
16) Anthracene	16.873	178	6481	0.343	ng/ul	100
19) Fluoranthene	18.819	202	8473	0.384	ng/ul#	92
20) Pyrene	19.186	202	8680	0.378	ng/ul#	90
21) Benzo(a)anthracene	20.967	228	7934	0.360	ng/ul	99
22) Chrysene	21.017	228	9039	0.408	ng/ul	100
24) Benzo(b)fluoranthene	22.466	252	8433	0.368	ng/ul	96
25) Benzo(k)fluoranthene	22.507	252	9901	0.419	ng/ul#	95
26) Benzo(a)pyrene	22.989	252	7791	0.372	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.121	276	10885	0.385	ng/ul#	86
28) Dibenzo(a,h)anthracene	25.141	278	8047	0.363	ng/ul#	95
29) Benzo(g,h,i)perylene	25.738	276	8293	0.372	ng/ul#	94

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

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