

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112224\
 Data File : BN035265.D
 Acq On : 23 Nov 2024 05:47
 Operator : RC/JU
 Sample : PB165115BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS115

Quant Time: Nov 23 06:10:01 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 : Fri Nov 22 12:45:23 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.318	152	4100	0.400	ng/ul	0.00
4) Naphthalene-d8	10.069	136	10422	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.981	164	5912	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.747	188	11370	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.988	240	7713	0.400	ng/ul	# 0.00
23) Perylene-d12	23.089	264	8574	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1943	0.519	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.670	152	5301	0.347	ng/ul	0.00
18) Fluoranthene-d10	18.795	212	9233	0.434	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.010	88	5357	1.317	ng/ul#	87
5) Naphthalene	10.113	128	9501	0.362	ng/ul	99
7) 2-Methylnaphthalene	11.747	142	6282	0.345	ng/ul	98
8) 1-Methylnaphthalene	11.972	142	6363	0.346	ng/ul	99
10) Acenaphthylene	13.690	152	9502	0.405	ng/ul	99
11) Acenaphthene	14.041	153	6886	0.384	ng/ul	98
12) Fluorene	15.041	166	7527	0.357	ng/ul	98
14) Pentachlorophenol	16.401	266	2048	0.773	ng/ul	98
15) Phenanthrene	16.785	178	11330	0.377	ng/ul	99
16) Anthracene	16.878	178	10171	0.365	ng/ul	99
19) Fluoranthene	18.828	202	11622	0.433	ng/ul#	95
20) Pyrene	19.195	202	11701	0.419	ng/ul#	91
21) Benzo(a)anthracene	20.973	228	10411	0.389	ng/ul	99
22) Chrysene	21.023	228	10806	0.401	ng/ul	99
24) Benzo(b)fluoranthene	22.472	252	11198	0.395	ng/ul	95
25) Benzo(k)fluoranthene	22.513	252	11747	0.402	ng/ul#	94
26) Benzo(a)pyrene	22.998	252	10195	0.394	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.138	276	14332	0.410	ng/ul#	84
28) Dibenzo(a,h)anthracene	25.151	278	10581	0.386	ng/ul#	92
29) Benzo(g,h,i)perylene	25.752	276	11119	0.403	ng/ul#	92

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

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