

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112224\
 Data File : BN035249.D
 Acq On : 22 Nov 2024 19:34
 Operator : RC/JU
 Sample : PB165200BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS200

Quant Time: Nov 22 22:32:11 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	7.322	152	3041	0.400	ng/ul	0.00
4) Naphthalene-d8	10.070	136	7749	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.981	164	4800	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.747	188	10585	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.988	240	7608	0.400	ng/ul	# 0.00
23) Perylene-d12	23.092	264	6707	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.976	96	1745	0.628	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.670	152	4104	0.361	ng/ul	0.00
18) Fluoranthene-d10	18.796	212	10030	0.478	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.006	88	4845	1.606	ng/ul#	90
5) Naphthalene	10.119	128	7130	0.365	ng/ul	100
7) 2-Methylnaphthalene	11.747	142	4869	0.360	ng/ul	99
8) 1-Methylnaphthalene	11.972	142	4930	0.361	ng/ul	98
10) Acenaphthylene	13.690	152	7708	0.405	ng/ul	99
11) Acenaphthene	14.041	153	5721	0.393	ng/ul	97
12) Fluorene	15.045	166	6521	0.381	ng/ul	99
14) Pentachlorophenol	16.405	266	2049	0.831	ng/ul	98
15) Phenanthrene	16.789	178	10867	0.388	ng/ul	100
16) Anthracene	16.878	178	9817	0.378	ng/ul	99
19) Fluoranthene	18.828	202	12651	0.478	ng/ul#	92
20) Pyrene	19.195	202	12901	0.468	ng/ul#	90
21) Benzo(a)anthracene	20.973	228	10486	0.397	ng/ul	99
22) Chrysene	21.026	228	11180	0.421	ng/ul	100
24) Benzo(b)fluoranthene	22.472	252	9560	0.431	ng/ul	95
25) Benzo(k)fluoranthene	22.516	252	10082	0.441	ng/ul#	95
26) Benzo(a)pyrene	23.001	252	8383	0.414	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.141	276	10775	0.394	ng/ul#	75
28) Dibenzo(a,h)anthracene	25.158	278	8000	0.373	ng/ul#	94
29) Benzo(g,h,i)perylene	25.758	276	8379	0.388	ng/ul#	92

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

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