

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112624\
 Data File : BN035331.D
 Acq On : 27 Nov 2024 00:54
 Operator : RC/JU
 Sample : PB165279BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS279

Quant Time: Nov 27 01:30:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112624.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	7.310	152	3271	0.400	ng/ul	0.00
4) Naphthalene-d8	10.059	136	8419	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.967	164	4912	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.734	188	9080	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.976	240	5734	0.400	ng/ul	# 0.00
23) Perylene-d12	23.071	264	6466	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.972	96	2086	0.698	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.659	152	4319	0.350	ng/ul	0.00
18) Fluoranthene-d10	18.786	212	6916	0.437	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.001	88	5515	1.699	ng/ul#	88
5) Naphthalene	10.108	128	7500	0.354	ng/ul	100
7) 2-Methylnaphthalene	11.736	142	5007	0.340	ng/ul	99
8) 1-Methylnaphthalene	11.961	142	5061	0.341	ng/ul	99
10) Acenaphthylene	13.681	152	7215	0.370	ng/ul	100
11) Acenaphthene	14.032	153	5516	0.370	ng/ul	99
12) Fluorene	15.031	166	5994	0.343	ng/ul	98
14) Pentachlorophenol	16.392	266	1381	0.653	ng/ul	98
15) Phenanthrene	16.776	178	8800	0.367	ng/ul	99
16) Anthracene	16.869	178	7585	0.340	ng/ul	100
19) Fluoranthene	18.814	202	8531	0.427	ng/ul#	93
20) Pyrene	19.186	202	8504	0.409	ng/ul#	93
21) Benzo(a)anthracene	20.961	228	6972	0.350	ng/ul	99
22) Chrysene	21.014	228	7845	0.392	ng/ul	100
24) Benzo(b)fluoranthene	22.454	252	7804	0.365	ng/ul	96
25) Benzo(k)fluoranthene	22.495	252	8433	0.383	ng/ul#	96
26) Benzo(a)pyrene	22.978	252	7023	0.360	ng/ul#	96
27) Indeno(1,2,3-cd)pyrene	25.111	276	9889	0.375	ng/ul#	87
28) Dibenzo(a,h)anthracene	25.128	278	7496	0.362	ng/ul#	94
29) Benzo(g,h,i)perylene	25.725	276	7833	0.377	ng/ul#	93

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

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