

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035575.D
 Acq On : 11 Dec 2024 20:18
 Operator : RC/JU
 Sample : PB165517BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS517

Quant Time: Dec 11 22:10:26 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 Dec 10 16:10:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.289	152	2446	0.400	ng/ul	0.00
4) Naphthalene-d8	10.037	136	6982	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	4752	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	10037	0.400	ng/ul	0.00
17) Chrysene-d12	20.961	240	7524	0.400	ng/ul	0.00
23) Perylene-d12	23.051	264	7293	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	1630	0.729	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.637	152	4177	0.408	ng/ul	0.00
18) Fluoranthene-d10	18.768	212	9203	0.443	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	4453	1.835	ng/ul	93
5) Naphthalene	10.081	128	6656	0.378	ng/ul	100
7) 2-Methylnaphthalene	11.714	142	4562	0.374	ng/ul	98
8) 1-Methylnaphthalene	11.939	142	4659	0.378	ng/ul	98
10) Acenaphthylene	13.662	152	7031	0.373	ng/ul	99
11) Acenaphthene	14.014	153	5439	0.377	ng/ul	98
12) Fluorene	15.013	166	6176	0.365	ng/ul	99
14) Pentachlorophenol	16.380	266	1946	0.832	ng/ul	98
15) Phenanthrene	16.760	178	10040	0.378	ng/ul	100
16) Anthracene	16.852	178	9019	0.366	ng/ul	100
19) Fluoranthene	18.800	202	10858	0.414	ng/ul	98
20) Pyrene	19.167	202	11374	0.417	ng/ul	96
21) Benzo(a)anthracene	20.947	228	9975	0.382	ng/ul	99
22) Chrysene	20.999	228	10507	0.400	ng/ul	100
24) Benzo(b)fluoranthene	22.440	252	9379	0.389	ng/ul	98
25) Benzo(k)fluoranthene	22.481	252	10456	0.421	ng/ul	97
26) Benzo(a)pyrene	22.963	252	8533	0.387	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.084	276	10031	0.337	ng/ul#	84
28) Dibenzo(a,h)anthracene	25.101	278	8001	0.343	ng/ul#	97
29) Benzo(g,h,i)perylene	25.698	276	7914	0.337	ng/ul	97

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

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