

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN121024\
 Data File : BN035547.D
 Acq On : 11 Dec 2024 00:30
 Operator : RC/JU
 Sample : PB165439BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS439

Quant Time: Dec 11 00:57:38 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	7.293	152	2685	0.400	ng/ul	0.00
4) Naphthalene-d8	10.036	136	7128	0.400	ng/ul	0.00
9) Acenaphthene-d10	13.949	164	4572	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.717	188	9034	0.400	ng/ul	# 0.00
17) Chrysene-d12	20.961	240	5892	0.400	ng/ul	# 0.00
23) Perylene-d12	23.051	264	6150	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.963	96	1810	0.738	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.642	152	3899	0.373	ng/ul	0.00
18) Fluoranthene-d10	18.767	212	7240	0.445	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.993	88	5045	1.894	ng/ul	90
5) Naphthalene	10.086	128	6723	0.374	ng/ul	99
7) 2-Methylnaphthalene	11.714	142	4410	0.354	ng/ul	97
8) 1-Methylnaphthalene	11.939	142	4607	0.366	ng/ul	99
10) Acenaphthylene	13.662	152	6653	0.367	ng/ul	99
11) Acenaphthene	14.014	153	5238	0.377	ng/ul	99
12) Fluorene	15.017	166	5786	0.355	ng/ul	99
14) Pentachlorophenol	16.379	266	1328	0.631	ng/ul	98
15) Phenanthrene	16.759	178	8999	0.377	ng/ul	100
16) Anthracene	16.848	178	7659	0.346	ng/ul	100
19) Fluoranthene	18.800	202	8950	0.436	ng/ul#	96
20) Pyrene	19.167	202	9302	0.436	ng/ul#	94
21) Benzo(a)anthracene	20.947	228	7264	0.355	ng/ul	98
22) Chrysene	20.999	228	8447	0.411	ng/ul	99
24) Benzo(b)fluoranthene	22.440	252	7802	0.384	ng/ul	96
25) Benzo(k)fluoranthene	22.481	252	8835	0.422	ng/ul#	97
26) Benzo(a)pyrene	22.960	252	7124	0.383	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.077	276	10435	0.416	ng/ul#	84
28) Dibenzo(a,h)anthracene	25.098	278	7211	0.367	ng/ul#	97
29) Benzo(g,h,i)perylene	25.695	276	7995	0.404	ng/ul#	95

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

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