

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035025.D
 Acq On : 12 Nov 2024 03:14
 Operator : RC/JU
 Sample : PB164722BS
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS722

Quant Time: Nov 12 03:43:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN110924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 09 00:09:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.562	152	3573	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.328	136	8789	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.199	164	5547	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	11583	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	10891	0.400	ng/ul	# 0.00
23) Perylene-d12	23.305	264	12412	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1866	0.549	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	4689	0.348	ng/ul	-0.01
18) Fluoranthene-d10	18.981	212	10481	0.375	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.182	88	4762	1.278	ng/ul#	92
5) Naphthalene	10.372	128	7001	0.318	ng/ul	98
7) 2-Methylnaphthalene	12.000	142	4907	0.299	ng/ul	100
8) 1-Methylnaphthalene	12.220	142	4922	0.299	ng/ul	99
10) Acenaphthylene	13.912	152	7195	0.317	ng/ul	99
11) Acenaphthene	14.259	153	5374	0.313	ng/ul	99
12) Fluorene	15.253	166	6134	0.281	ng/ul	99
14) Pentachlorophenol	16.603	266	1955	0.730	ng/ul	97
15) Phenanthrene	16.988	178	9853	0.318	ng/ul	99
16) Anthracene	17.076	178	9249	0.318	ng/ul	99
19) Fluoranthene	19.009	202	12105	0.333	ng/ul#	95
20) Pyrene	19.372	202	12600	0.334	ng/ul#	96
21) Benzo(a)anthracene	21.128	228	13165	0.334	ng/ul	99
22) Chrysene	21.177	228	13039	0.324	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	13671	0.313	ng/ul	95
25) Benzo(k)fluoranthene	22.703	252	14109	0.309	ng/ul#	96
26) Benzo(a)pyrene	23.212	252	12899	0.331	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.457	276	17074	0.358	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.473	278	13328	0.359	ng/ul#	97
29) Benzo(g,h,i)perylene	26.107	276	13339	0.366	ng/ul#	97

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

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