

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035005.D
 Acq On : 11 Nov 2024 15:13
 Operator : RC/JU
 Sample : PB164747BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS747

Quant Time: Nov 11 16:04:24 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	3101	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.328	136	7731	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.199	164	5020	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	10784	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	10049	0.400	ng/ul	# 0.00
23) Perylene-d12	23.308	264	10758	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	2443	0.828	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	6117	0.517	ng/ul	-0.01
18) Fluoranthene-d10	18.981	212	13946	0.540	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.182	88	6887	2.130	ng/ul#	91
5) Naphthalene	10.377	128	9720	0.502	ng/ul	98
7) 2-Methylnaphthalene	12.000	142	6851	0.474	ng/ul	100
8) 1-Methylnaphthalene	12.220	142	7996	0.551	ng/ul	99
10) Acenaphthylene	13.916	152	9897	0.483	ng/ul	99
11) Acenaphthene	14.263	153	7498	0.482	ng/ul	99
12) Fluorene	15.253	166	8588	0.434	ng/ul	98
14) Pentachlorophenol	16.599	266	2397	0.961	ng/ul	98
15) Phenanthrene	16.987	178	13921	0.483	ng/ul	98
16) Anthracene	17.080	178	13043	0.482	ng/ul	98
19) Fluoranthene	19.009	202	16947	0.505	ng/ul#	96
20) Pyrene	19.372	202	17497	0.502	ng/ul#	96
21) Benzo(a)anthracene	21.128	228	18013	0.495	ng/ul	98
22) Chrysene	21.177	228	18125	0.488	ng/ul	99
24) Benzo(b)fluoranthene	22.665	252	18685	0.494	ng/ul	98
25) Benzo(k)fluoranthene	22.706	252	19058	0.481	ng/ul	98
26) Benzo(a)pyrene	23.211	252	16736	0.495	ng/ul#	98
27) Indeno(1,2,3-cd)pyrene	25.460	276	22294	0.539	ng/ul	94
28) Dibenzo(a,h)anthracene	25.477	278	17686	0.549	ng/ul	98
29) Benzo(g,h,i)perylene	26.114	276	17372	0.550	ng/ul	98

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

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