

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111124\
 Data File : BN035048.D
 Acq On : 12 Nov 2024 22:24
 Operator : RC/JU
 Sample : PB164733BS
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS733

Quant Time: Nov 12 22:54:52 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	7.558	152	3525	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.322	136	8605	0.400	ng/ul	#-0.02
9) Acenaphthene-d10	14.194	164	5706	0.400	ng/ul	-0.01
13) Phenanthrene-d10	16.945	188	12603	0.400	ng/ul	0.00
17) Chrysene-d12	21.136	240	11835	0.400	ng/ul	#-0.01
23) Perylene-d12	23.302	264	12615	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	1907	0.569	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.917	152	4513	0.342	ng/ul	-0.02
18) Fluoranthene-d10	18.977	212	11295	0.372	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.182	88	4771	1.298	ng/ul#	92
5) Naphthalene	10.372	128	6800	0.316	ng/ul	97
7) 2-Methylnaphthalene	11.994	142	4811	0.299	ng/ul	99
8) 1-Methylnaphthalene	12.214	142	4826	0.299	ng/ul	98
10) Acenaphthylene	13.912	152	7095	0.304	ng/ul	100
11) Acenaphthene	14.259	153	5473	0.310	ng/ul	98
12) Fluorene	15.249	166	6407	0.285	ng/ul	99
14) Pentachlorophenol	16.599	266	1697	0.582	ng/ul	98
15) Phenanthrene	16.983	178	10672	0.317	ng/ul	99
16) Anthracene	17.076	178	9850	0.311	ng/ul	99
19) Fluoranthene	19.004	202	13156	0.333	ng/ul#	95
20) Pyrene	19.367	202	13707	0.334	ng/ul#	95
21) Benzo(a)anthracene	21.122	228	14029	0.327	ng/ul	100
22) Chrysene	21.174	228	14265	0.326	ng/ul	99
24) Benzo(b)fluoranthene	22.656	252	14730	0.332	ng/ul	98
25) Benzo(k)fluoranthene	22.700	252	14986	0.323	ng/ul#	99
26) Benzo(a)pyrene	23.206	252	13169	0.332	ng/ul#	98
27) Indeno(1,2,3-cd)pyrene	25.450	276	17741	0.366	ng/ul#	93
28) Dibenzo(a,h)anthracene	25.463	278	13898	0.368	ng/ul#	98
29) Benzo(g,h,i)perylene	26.100	276	13942	0.377	ng/ul	98

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

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