

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034990.D
 Acq On : 10 Nov 2024 21:41
 Operator : RC/JU
 Sample : PB164741BS
 Misc :
 ALS Vial : 84 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS741

Quant Time: Nov 10 22:28:59 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	7.566	152	2259	0.400	ng/ul	0.00
4) Naphthalene-d8	10.328	136	5479	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.199	164	4033	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.945	188	9849	0.400	ng/ul	0.00
17) Chrysene-d12	21.140	240	10165	0.400	ng/ul	0.00
23) Perylene-d12	23.302	264	10912	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.153	96	1874	0.872	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.923	152	3848	0.459	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	12082	0.463	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.183	88	4160	1.766	ng/ul#	91
5) Naphthalene	10.378	128	5313	0.388	ng/ul	99
7) 2-Methylnaphthalene	12.000	142	3829	0.374	ng/ul	98
8) 1-Methylnaphthalene	12.220	142	3885	0.378	ng/ul	99
10) Acenaphthylene	13.917	152	6185	0.375	ng/ul	99
11) Acenaphthene	14.264	153	4722	0.378	ng/ul	99
12) Fluorene	15.254	166	5838	0.367	ng/ul	99
14) Pentachlorophenol	16.599	266	1796	0.789	ng/ul	97
15) Phenanthrene	16.988	178	10248	0.389	ng/ul	99
16) Anthracene	17.076	178	9583	0.387	ng/ul	99
19) Fluoranthene	19.009	202	13318	0.392	ng/ul#	97
20) Pyrene	19.372	202	14024	0.398	ng/ul#	97
21) Benzo(a)anthracene	21.125	228	14266	0.388	ng/ul	99
22) Chrysene	21.178	228	14734	0.392	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	14926	0.389	ng/ul	99
25) Benzo(k)fluoranthene	22.703	252	16070	0.400	ng/ul	99
26) Benzo(a)pyrene	23.209	252	13824	0.403	ng/ul#	99
27) Indeno(1,2,3-cd)pyrene	25.457	276	18727	0.447	ng/ul#	94
28) Dibenzo(a,h)anthracene	25.470	278	14638	0.448	ng/ul	98
29) Benzo(g,h,i)perylene	26.107	276	14478	0.452	ng/ul	97

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

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