

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110924\
 Data File : BN034960.D
 Acq On : 10 Nov 2024 02:36
 Operator : RC/JU
 Sample : PB164807BS
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS807

Quant Time: Nov 10 03:00:43 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	2629	0.400	ng/ul	0.00
4) Naphthalene-d8	10.333	136	6923	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.203	164	5588	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.949	188	13653	0.400	ng/ul	0.00
17) Chrysene-d12	21.142	240	13367	0.400	ng/ul	0.00
23) Perylene-d12	23.308	264	12160	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.149	96	1803	0.721	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.928	152	5085	0.480	ng/ul	0.00
18) Fluoranthene-d10	18.981	212	16691	0.486	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.182	88	4914	1.792	ng/ul#	92
5) Naphthalene	10.383	128	7122	0.411	ng/ul	98
7) 2-Methylnaphthalene	12.005	142	5410	0.418	ng/ul	99
8) 1-Methylnaphthalene	12.225	142	6308	0.486	ng/ul	99
10) Acenaphthylene	13.921	152	9056	0.397	ng/ul	99
11) Acenaphthene	14.268	153	6866	0.397	ng/ul	99
12) Fluorene	15.258	166	8513	0.387	ng/ul	99
14) Pentachlorophenol	16.603	266	2502	0.793	ng/ul	98
15) Phenanthrene	16.992	178	14885	0.408	ng/ul	99
16) Anthracene	17.080	178	13648	0.398	ng/ul	99
19) Fluoranthene	19.014	202	19043	0.426	ng/ul#	97
20) Pyrene	19.376	202	19866	0.429	ng/ul	98
21) Benzo(a)anthracene	21.125	228	19298	0.399	ng/ul	99
22) Chrysene	21.177	228	19787	0.401	ng/ul	100
24) Benzo(b)fluoranthene	22.662	252	18156	0.425	ng/ul	99
25) Benzo(k)fluoranthene	22.706	252	18393	0.411	ng/ul	100
26) Benzo(a)pyrene	23.211	252	15385	0.403	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.460	276	18221	0.390	ng/ul#	95
28) Dibenzo(a,h)anthracene	25.473	278	14396	0.396	ng/ul	98
29) Benzo(g,h,i)perylene	26.114	276	14080	0.395	ng/ul	98

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

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